

SCREENING SITE INSPECTION REPORT  
FOR  
HADER GROUNDWATER CONTAMINATION  
HADER, MINNESOTA

U.S. EPA ID: MND981961873

EPA Region 5 Records Ctr.



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## EXECUTIVE SUMMARY

On March 29, 1988, the Minnesota Pollution Control Agency (MPCA) staff conducted a Screening Site Investigation (SSI) at the Hader Ground Water Contamination Site (Site), Goodhue County, Minnesota. The purpose of the SSI was to determine if high amounts of lead, nickel, sulfate, and iron in the ground water could be attributed to a contaminant source.

Ten residential wells were sampled in August 1987, and three were found to exceed the Recommended Allowable Limits (RALs) established by the Minnesota Department of Health (MDH) for lead and nickel and U.S. Environmental Protection Agency (EPA) Secondary Drinking Water Standards for iron and sulfate. Upon receipt of these results the MPCA began supplying these households with bottled drinking water.

After further screening of households that had reported poor quality water, seven households were chosen and sampled in April 1988 for volatile organic aromatics (VOAs), Acid Base Neutrals (ABNs) compounds, pesticides, polychlorinated biphenyls (PCBs), dissolved metals, nitrates, nitrites, sulfates, and chlorides. These compounds and analytes are on the U.S. EPA Target Compound List (TCL) and Target Analyte List (TAL). Organic compounds were all below the detection limits of laboratory equipment. Elevated levels of inorganics included iron, zinc, sulfate and nickel. Three of the seven households sampled had concentrations of nickel above RALS.

## 1.0 INTRODUCTION

The MPCA, working under a Cooperative Agreement with the U.S. Environmental Protection Agency (EPA), has conducted a SSI at the Hader Ground Water Contamination Site, Minneola Township, Goodhue County.

The Site was initially discovered as a result of a residential complaint of extremely poor quality ground water, i.e., cloudy, poor taste, and high iron content.

The Site was placed on the U.S. EPA's Comprehensive Environmental Response, Compensation, and Liability Information Systems (CERCLIS) list inventory on July 30, 1987. A Preliminary Assessment (PA) of the Site was conducted by MPCA staff and submitted to U.S. EPA on September 30, 1987. The MPCA staff prepared and submitted an SSI workplan for review by the U.S. EPA Region V staff. The MPCA staff received approval from U.S. EPA for the SSI workplan in January 1988.

The objectives of an SSI are stated in a directive from U.S. EPA which outlines Pre-remedial Strategies under the Comprehensive Environmental Response, Compensation and Liability Act (CERCLA). The directive states:

All sites will receive an SSI in an effort to; 1) collect additional data beyond the scope of the PA to enable a more refined preliminary Hazardous Ranking System (HRS) score, 2) establish priorities among sites most likely to qualify for the National Priorities List (NPL), and 3) identify the data requirements for a Listing Site Investigation (LSI).

An SSI will not have rigorous data quality objectives (DOQs). Based on the refined preliminary HRS score and other technical judgement factors, the site will then either be designated as NFRAP [no further remedial action planned], or carried forward as an NPL listing candidate. A listing SI will not automatically be done on these sites, however. First, they will go through a management evaluation to determine whether they can be addressed by another authority such as, RCRA [Resource Conservation and Recovery Act] . . . Sites that are designated NFRAP or deferred to other statutes are not candidates for a listing SI.

The listing SI will address all the data requirements of the revised HRS using field screening and NPL level DOQs. It may also provide needed data in a format to support remedial investigation workplan development. Only sites that appear to score high enough for listing and that have not been deferred to another authority will receive a listing SI (9).

U.S. EPA Region V has also instructed the MPCA to identify sites during the SSI that may require prompt removal action to remediate an immediate human health and/or environmental threat.

## 2.0 SITE INFORMATION

### 2.1 Site Description

The Site is located in a sparsely populated rural agricultural area of southeastern Minnesota, in which dairy farming is the predominant industry. The boundaries of the Site are non-distinct and cover a general area of

HADER GROUNDWATER CONTAMINATION  
HADER, MN

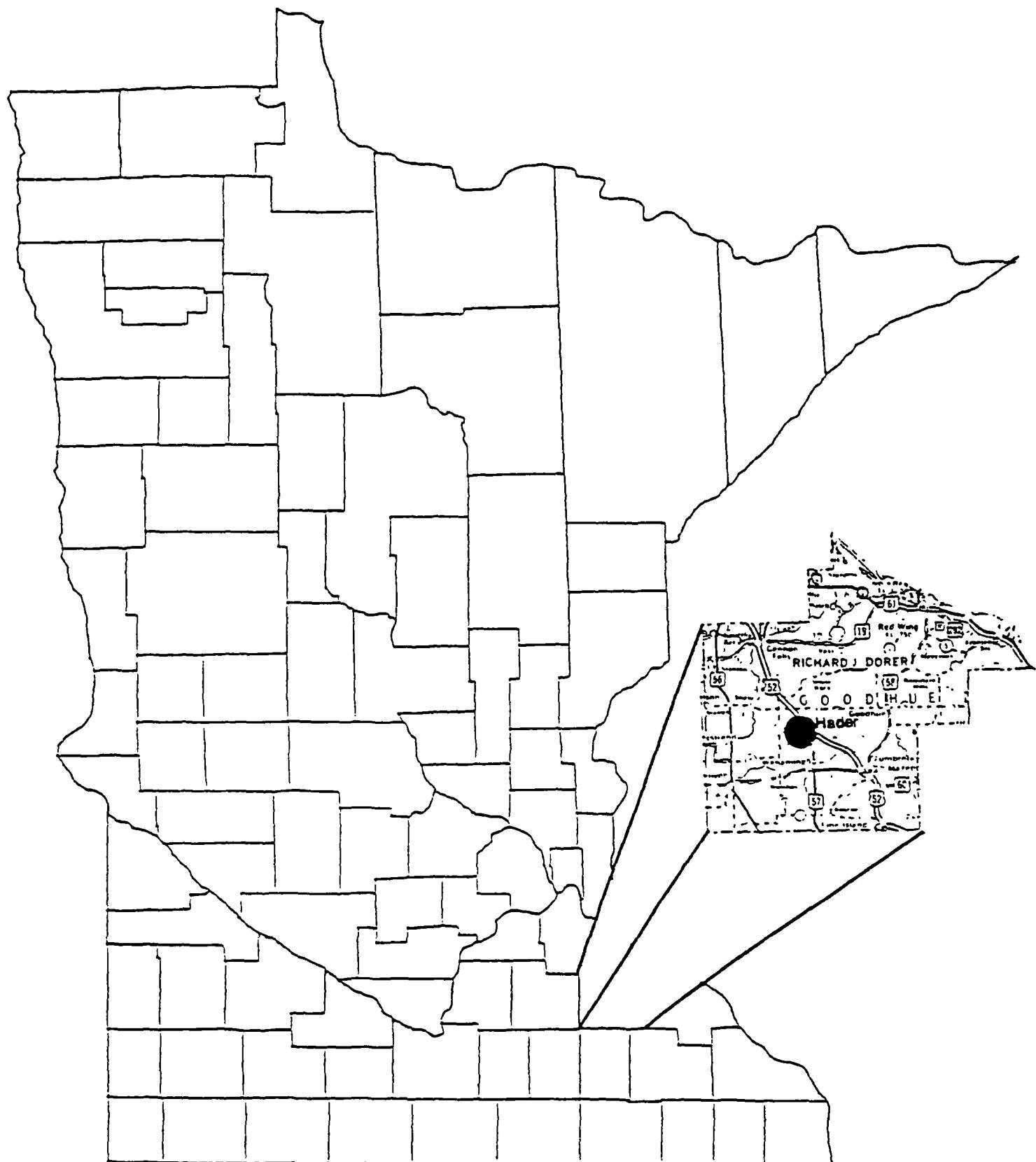


Figure I  
SITE LOCATION MAP

approximately 24 square miles (Figure I and Appendix A). Belle Creek and Minneaola townships comprise the majority of the study area. The nearest population center is Zumbrota, approximately 8 miles southeast of the Site.

## 2.2 History

The area was first brought to the attention of the MPCA through a complaint filed by a resident in the area. The complainant reported cloudy residues and high iron content in the drinking water, which readily formed a film or stain upon contact with piping, containers, human skin, and clothing. Both human and livestock health problems were reported if the water was used for any length of time for drinking purposes. Further interviews with other residents in the area resulted in similar reports of poor quality drinking water, and health problems. Although no medical evidence confirms that the ground water at the Site is a cause of health problems, there appears to be a relationship between ground water consumption and poor health of some of the residents and their livestock.

In August 1987, MPCA staff conducted water sampling at ten homes in the area, including that of the complainant. A limited number of metals, volatile hydrocarbons, and pesticides were analyzed for by the Minnesota Department of Health (MDH) laboratories in Minneapolis, Minnesota. Of the ten samples, three were found to exceed MDH RALS for lead and nickel and U.S. EPA Secondary Drinking Water Standards for iron and sulfate. MPCA staff also noted that wells, in which elevated levels of metals were measured, also had extremely low pH levels. Difficulty was encountered in stabilizing the well prior to sampling, as pH would often continue to drop after numerous well volumes were extracted. A pH as low as 4.8 was recorded at the complainants well [pH

levels of this magnitude do not usually occur in ground water under natural conditions].

Upon receipt of these results, the MPCA notified the residents and MDH of the poor quality of their drinking water and began supplying bottled water to the three homes which exceeded drinking water standards.

### 3.0 SSI OBJECTIVES

The principal objective of the SSI was to determine if the elevated levels of inorganic compounds detected in domestic wells in the vicinity of Minneola and Belle Creek Townships, Minnesota are attributable to a contaminant source or sources. If a contaminant source was identified, and an observed release to the ground or surface water was documented, further investigation would be warranted. If a contaminant source was not identified, the Site would receive a "No Further Remedial Action Planned" (NFRAP) priority.

### 4.0 GEOLOGY

#### 4.1 Physical and Hydrogeologic Properties

The Site is underlain by Cambrian, Ordovician, Cretaceous bedrock, and Quaternary glacial and surficial deposits (Figure II). The regional bedrock slopes toward the southeast, while the regional ground water flow is to the north, toward the Cannon River (10).

The oldest geologic formation underlying the Site is the Pre - Cambrian bedrock system of igneous and sedimentary rocks (Figure II). The oldest

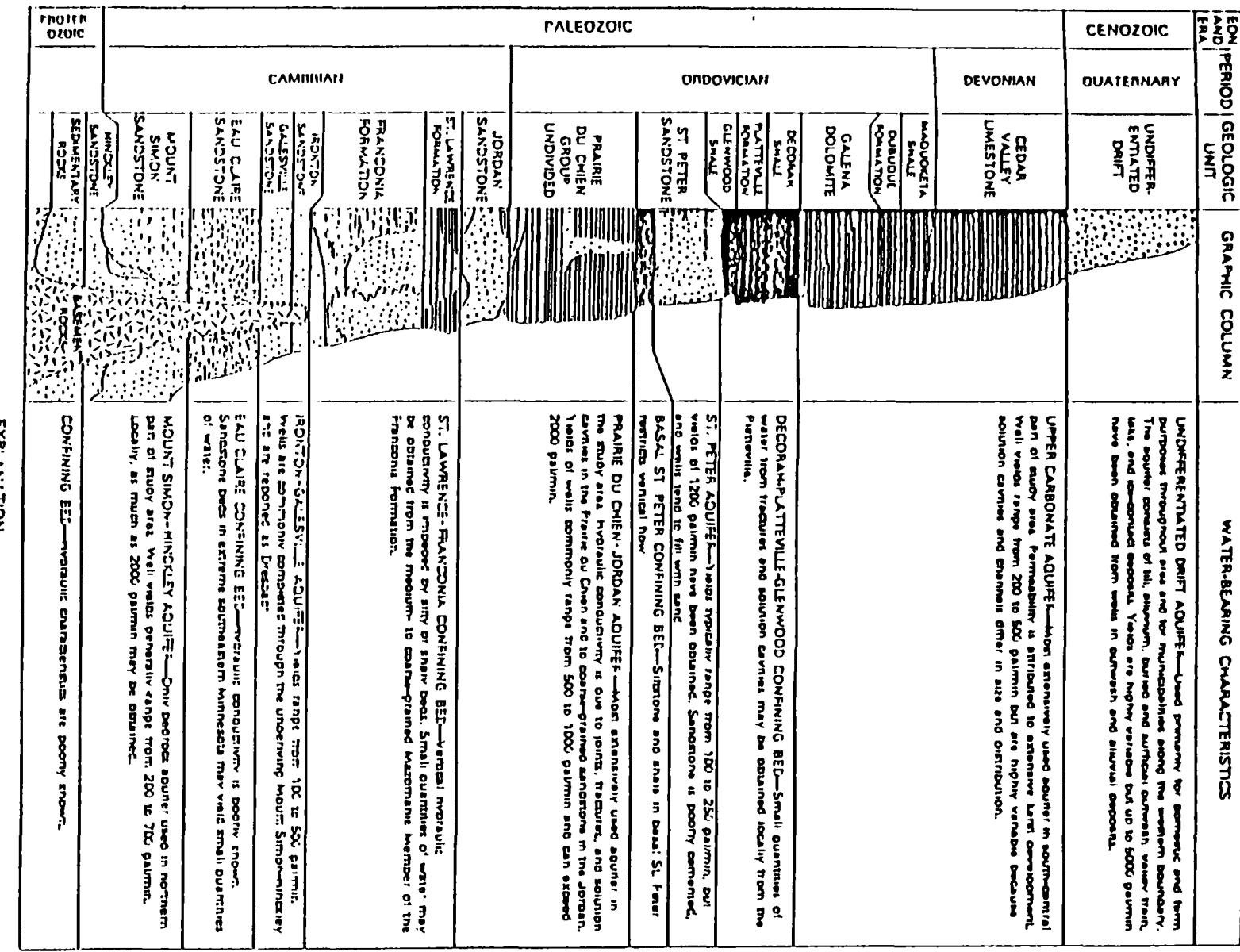


Figure 11

## AQUIFER UNITS AND PRODUCTIVITY

sedimentary bedrock is the Fond du Lac formation composed of well cemented, fine-grained arkosic sand to siltstone with interbedded shales. Tertiary igneous intrusions extend vertically through Pre-Cambrian sedimentary basement rocks into the Ironton formation. The Fond du Lac and areas of igneous intrusions are not utilized for ground water supplies in southeastern Minnesota.

The deepest primary water bearing units are the Hinkley and Mt. Simon formations (Figure II). The Hinkley formation is a medium to coarse grained arkosic sandstone which grades into the overlying Mt. Simon formation. The Mt. Simon formation is a fine to coarse grained quartzose with thin shale beds in the upper portion. Due to the absence of a confining layer between the Mt. Simon and Hinkley formations, they are considered to be one hydrologic unit. Water yields are moderate to high depending upon placement of wells. Hydraulic conductivity varies from 2 to 23 feet/day (5).

The overlying Eau Claire formation acts as a confining layer between the Mt. Simon-Hinkley aquifer and the Galesville-Ironton formations (Figure II). As with the Mt. Simon-Hinkley formations, there is the absence of a confining layer between the Galesville and Ironton formations. The lack of a confining layer allows the two sandstones to act as one hydraulic unit. The Ironton-Galesville Group is a medium to fine-grained sandstone. Average yields range from 4 to 33 feet/day (5).

The Franconia formation is a fine-grained sandstone with shale layers interspersed in some areas (Figure II). Due to the interfingering of the shale layers, the hydraulic potential of the Franconia sandstone is limited to the upper Reno unit which has less shale. Over all the formation acts as a

confining unit (5).

The St. Lawrence formation is a dolomitic shale with areas of sandstone and greensand (Figure II). The St. Lawrence is approximately five feet thick, and with the Franconia formation acts as a confining layer between the underlying Franconia and the overlying Jordan sandstone (5).

The Jordan sandstone is massive with numerous areas of cross bedding and an average thickness of 100 feet (Figure II). Pyrite stringers are interspersed throughout the formation and hematite cementing is common. Due to the absence of a confining layer, the Jordan acts as a primary aquifer with the interconnected Prairie du Chien formation, providing large yields of acceptable quality drinking water in the upland regions (7).

The Prairie du Chien formation of Ordovician age is comprised of two units, the Oneata and Shakopee dolomites (Figure II). The Oneata unit averages 150 feet in thickness where it outcrops along the Cannon River, north of the Site (10). The Oneata dolomite has limited hydraulic potential yields along joints and bedding planes, whereas the overlying Shakopee dolomite has a higher potential yield due to interspersed sandy layers.

As previously stated, there is no confining unit between the Prairie du Chien aquifer and the Jordan aquifer, therefore ground water mixing between the two formations is possible. Of the two units (Prairie du Chien and Jordan), the Jordan sandstone has greater porosity but the Prairie du Chien Group has a higher fracture permeability and therefore greater yields. The average hydraulic conductivity for the Prairie du Chien - Jordan aquifer is 5 to 67 feet/day (5).

The St. Peter formation of Ordovician age thins in the western margin of Goodhue County, relative to the eastern margin. The average thickness of the St. Peter underlying the Site is 75 feet. The St. Peter formation is a well sorted, clean sandstone with pyrite stringers often interspersed (Figure II). Water resources are adequate for domestic use, with average hydraulic conductivity of 3 to 33 feet/day, but yields are restricted in areas where the formation has been eroded to a thickness as little as 10 feet (5).

The Glenwood formation is thin (3 feet) to non-existent in the vicinity of the Site. The formation is a fissile shale and when present acts as a confining layer between the underlying St. Peter sandstone and the overlying Platteville limestone (5).

Unlike the Glenwood formation, the Platteville formation is laterally extensive across the Site with an average thickness of 25 feet. Fractures and joints in the dolomitic limestone provide adequate porosity and permeability for limited domestic and livestock yields (7).

Overlying the Platteville formation is the Decorah shale (Figure II). The Decorah formation is thin (less than 5 feet) to non-existent at the Site. When present, the shale is greenish-gray in color and ranges from fissile to blocky and acts as a confining layer between the underlying Platteville limestone and the Galena dolomite. When absent, there is direct hydraulic connection between the Platteville formation and the Galena dolomite. Due to the interfingering of shales within the Platteville formation, the Decorah - Platteville - Glenwood formations are often viewed as a confining layer on a regional scale (5).

The Galena formation overlies the Decorah Shale at the Site, unless removed by surface erosion (Figure II). The Galena formation is a dolomite exhibiting karst characteristics throughout southeastern Minnesota. Solution tunnels, caves, and hummocky topography are primary indicators for the presence of the Galena dolomite. The Galena dolomite has been quarried throughout Goodhue County for use in concrete and other building purposes. Several abandoned quarries in the Galena dolomite are within the 24 squares miles established as the Site. The Galena formation is not widely utilized for domestic ground water in the area due to its unreliability for ground water yields in the absence of overlying formations. The solutioned nature of the Galena also makes the formation susceptible to ground water contamination (5 and 7).

Glacial deposits overlie the bedrock throughout the Site, unless absent due to erosion. The composition varies from sand and gravel outwash to clay tills. Ground water yields are adequate for domestic yields (<50 gal/day) with water quality somewhat variable. Due to recharge from precipitation, the water table is susceptible to seasonal variations (9).

#### 4.2 Residential Wells

Residential wells at the Site range in depth from 45 feet to 390 feet, therefore several aquifers were sampled (Appendix H). The glacial drift aquifer is the least used ground water source due to its susceptibility to fluctuating water table levels and contamination from pesticides and fertilizers. The Platteville limestone is used to a limited extent by households at the Site, but wells set in the Platteville aquifer have not reported any problems with the quality of the ground water.

The majority of the wells at the Site draw ground water from the lower part of the Prairie du Chien formation or the Shakopee limestone. One well is set at a greater depth in the Jordan sandstone (Appendix H). The Jordan aquifer is the deepest aquifer used by residents at the Site. The deeper Mt. Simon-Hinkley and Ironton-Galesville sandstone aquifers are not used at the Site. The towns of Wanamingo and Zumbrota, south of the Site, utilize the Mt. Simon-Hinkley formation for municipal ground water supplies (1).

All wells at the Site were constructed without the installation of a screen and often without grouting (this is a common practice for older water supply wells). Therefore, the water drawn from a well may be derived from a combination of aquifer sources, not solely from the aquifer in which the casing is terminated. This creates difficulty in determining the unit(s) from which a ground water sample was taken.

#### 4.3 Water Quality

The ground water in the vicinity of the Site is classified as a "calcium-magnesium bicarbonate type" (9). As a result, the most common non-contaminant related complaint concerns the hardness of the water. Some residents at the Site have corrected hardness problems with the installation of water softening devices.

A second common problem associated with ground water at the Site has been relatively high amounts of dissolved iron, as well as iron bacteria. The maximum recorded ambient iron level for the Prairie du Chien aquifer in Goodhue County is 620 ug/l (1 and Appendix D). The Jordan sandstone aquifer

is often plagued by high dissolved iron contents which range from 0.2 to .68 mg/l (5 and 9). When elevated levels of iron are reached, the ground water is often unusable from an aesthetic point due to poor taste and the water's tendency to stain clothing and plumbing. Iron bacteria may also induce staining and odor problems, as well.

## 5.0 SURFACE WATER

The Site is located within the Cannon River Watershed District. Belle Creek, the predominant surface water body, flows north through the Site to the Cannon River (9). In turn, local intermittent streams flow into Belle Creek. These intermittent streams are primarily used for livestock. Recreational use is limited due to the small size of the streams. Farm ponds are spread throughout the Site area and are used for watering livestock, recreational fishing, and swimming.

## 6.0 FIELD PROCEDURE

### 6.1 On-Site Interviews

Numerous interviews were taken with residents at the Site throughout the on-going SSI (MPCA files). Most residents could provide no evidence for suspected releases of hazardous substances to the environment in the immediate vicinity. An unpermitted dump in an abandoned quarry had been operating from the early 1970s through 1980, approximately 1 and a half miles to the northwest of the area established as the Site. The dump is being addressed under a separate SSI entitled "Hader Dump".

Some area residents observed the quality of their water also decreased when blasting began at a second separate quarry in the area. On days following a great deal of blasting, residents noticed cloudy residues and an increase in the amount of iron in their drinking water. Blasting has ceased at the second quarry.

It is also a common practice for residents without water treatment systems to chlorinate their wells in an effort to control iron bacteria, both which degrade the quality of the drinking water. The added shock of chlorine may also contribute to adverse health effects (3).

## 6.2 Sampling Procedure

Due to the large area established as the Site (approximately 24 square miles) and no direct evidence indicating possible contaminant sources, monitoring wells and soil borings were deemed inappropriate. Residential wells were selected on the basis of residential complaints and low pH levels.

Using a low pH as a criteria for selecting wells for sampling was justified by the correlation between low pH and high metal levels detected in previous non-contract lab testing performed in August 1987, by MPCA staff, through the MDH Environmental Laboratory in Minneapolis, Minnesota (Appendix C).

On April 29 through 31, 1988, seven residential wells were sampled for U.S. EPA TAL and TCL analytes and compounds including; VOAs, ABNs, pesticides/PCBs, metals, sulfates, chloride and nitrate/nitrite (Table I, Table II, Appendix C and Figure III). The original complainant has since moved from the area, therefore no sample was obtained due to restricted access to the well head. Special Analytical Services (SAS), as detailed in the Quality Assurance

non responsive

TABLE I  
HADER GROUND WATER CONTAMINATION  
RESIDENTIAL WELLS SAMPLED

<u>SAMPLE NUMBER</u>	<u>NAME AND ADDRESS</u>
S01	non responsive
S02	
S03	
S04	
S05, D05	
S06	
S07	
S08	

Table II  
 Summary of  
 Chemical Analysis for  
 Hader Ground Water Contamination

Sample Collection Information and Detected Parameters	<u>Sample Number</u>					
	S01	S02	S03	S04	S05	S06
Date	3-30-88	3-30-88	3-29-88	3-29-88	3-31-88	3-31-88
Time	12:31	10:33	11:53	14:49	11:36	11:56
Inorganic Traffic Report Number	MEW615	MEW616	MEW617	MEW618	MEW619	MEW620
Compound Detected (ug/l ppb unless indicated)						
Aluminum	633	----	25.9	1660	34.0	34.3B
Arsenic	1.7	---	---	---	---	8.4
Barium	34.4	88.2	72.6	91.5	---	28.8B
Calcium	167000	150000	74300	80000	111000	113000
Chromium	----	----	----	----	----	11.0
Cobalt	16.8	9.0	----	47.6	----	----
Copper	27.4	----	----	----	----	22.0
Iron	78800	5790	1560	83200	9780	9700
Lead	5.5	3.4	3.8	9.5	2.0	3.4
Magnesium	17100	40200	23200	20700	40600	41000
Manganese	570	181	68.8	655	621	574
Nickel	86.2	42.5	----	197	----	434

Table II

Summary of Chemical Analysis for  
Hader Ground Water Contamination

Sample Collection Information and Detected Parameters	Sample Number					
	S01	S02	S03	S04	S05	S06
Date:	3-30-88	3-30-88	3-29-88	3-29-88	3-31-88	3-31-88
Time:	12:31	10:33	11:53	14:49	11:36	11:56
Inorganic Traffic Report Number	MEW615	MEW616	MEW617	MEW618	MEW619	MEW620
Compound Detected (ug/L)						
Potassium	1320	1300	1030	833	17808	17808
Sodium	3570	2290	3420	1800	3230	2890
Vanadium	----	----	----	----	----	25.8
Zinc	1340	908	39.1	1470	32.3	28.5
Sulfate (mg/l ppm)	605	310	19.0	315	136	139
						890

Table II

Summary of Chemical Analysis for  
Hader Ground Water Contamination

Sample Collection Information and Detected Parameters	S07	S08	Sample Number					
Date:	3-31-88	3-31-88						
Time:	14:30	15:53						
Inorganic Traffic Report Number	MEW622	MEW623						
<u>Compound Detected (ug/L)</u>								
Aluminum	----	216						
Arsenic	3.6	5.0						
Barium	85.1	90.6						
Calcium	84000	85500						
Chromium	----	----						
Cobalt	----	120						
Copper	----	----						
Iron	3560	76800						
Lead	----	----						
Magnesium	28500	27000						
Manganese	69.8	517						

Table II

## **Summary of Chemical Analysis for Hader Ground Water Contamination**

Sample Collection Information and Detected Parameters	S07	S08	Sample Number				
			_____	_____	_____	_____	_____
Date:	3-31-88	3-31-88	_____	_____	_____	_____	_____
Time:	14:30	15:53	_____	_____	_____	_____	_____
Inorganic Traffic Report Number	MEW622	MEW623	_____	_____	_____	_____	_____
<u>Compound Detected (ug/L)</u>							
Nickel	---	342	_____	_____	_____	_____	_____
Potassium	5660	1160	_____	_____	_____	_____	_____
Sodium	6680	3600	_____	_____	_____	_____	_____
Vanadium	----	----	_____	_____	_____	_____	_____
Zinc	2720	477	_____	_____	_____	_____	_____
Sulfate (mg/l ppm)	24	314	_____	_____	_____	_____	_____

Project Plan (QAPP), were used to gain lower detection limits, given the samples were taken from drinking water wells (8). Photographs were taken to document each sampling point (Appendix B).

## 7.0 ANALYTICAL RESULTS

Results for the organic compounds showed no contaminants present above detection limits for all seven households sampled. Inorganic results showed the water to be a calcium-magnesium bicarbonate type, as defined by ambient data, that is extremely hard (Table I and Appendix E). High levels of iron were detected but these were not above the Secondary Drinking Water Standard of 500 ug/l. Sulfate levels were elevated at 15 households, but as with iron, they did not exceed the Secondary Drinking Water Standard of 300 ug/l. Lead was also detected, but was within the MDH RALs of 20 ug/l. Zinc exceeded EPAs Secondary Drinking Water Standard of 500 ug/l in 5 of the wells sampled. Nickel was found to be above the RAL of 150 ug/l at three of the households sampled at concentrations of 197, 434 and 342 ug/l.

Holding times were exceeded for extractions from semi-volatile and pesticide/PCBs fractions for all samples, therefore these results have been flagged as estimated (J=Positive, U=Negative) (Appendix C). Common laboratory contaminants were found in the laboratory blanks, but were all within acceptable limits. Further explanation of the data qualifiers preceeds the U.S. EPA Contract Laboratory Program (CLP) data in Appendix C.

## 8.0 CONCLUSIONS

Based on the data collected prior to and during the initiation of the SSI,

there appears that no hazardous substances have entered into the ground water system at the Site. Several natural explanations may account for the anomalously low pH and high metal levels.

The high amounts of iron may be attributed to three sources, pyrite deposits and interstitial hematite cement within the Prairie du Chien - Jordan aquifer and/or iron-precipitating bacteria multiplying within the well casing and the aquifer in which the well is located.

Hematite cement is a fairly common cement in sandstone and dissolves fairly easily, thus contributing to dissolved iron in the ground water. Pyrite deposits have been found in limited areas throughout the sedimentary rocks in southeastern Minnesota. The physical and chemical characteristics of the ground water in the vicinity of the Site may encourage the dissolution of pyrite, thus creating an increase in dissolved iron and sulfate in the ground water.

The third possible source of iron in the ground water is iron-precipitating bacteria. Iron-precipitating bacteria can cause a number of problems in water wells. The primary problem is water degradation through an increase of iron in the ground water (3). A secondary problem is encrustation and corrosion by the bacteria, resulting in poor well yields and degradation of the casing and pump (2). Some iron-precipitating bacteria can also lower the pH of the water in the well and surrounding aquifer (3). A low pH can result in the leaching of metals from well casings, pumps, the column pipe from the pump to the surface, and plumbing.

The MPCA staff feels that iron-precipitating bacteria may be contributing to

the conditions at the Site, allowing elevated metal levels to be detected. It is known that many residents at the Site and in the surrounding area experience iron-precipitating bacteria problems. Should the type of bacteria in the wells and aquifers lend itself to lowering the pH of the ground water, the leaching process may be extracting metals from various parts of the well system.

Specifically, lead may be extracted from lead plumbing pipes and spigots, while nickel can be leached from the brass fittings on submersible pumps. The coating applied to galvanized steel pump column pipes could be a source of zinc. Iron may be generated by the bacteria, which initiated the problem.

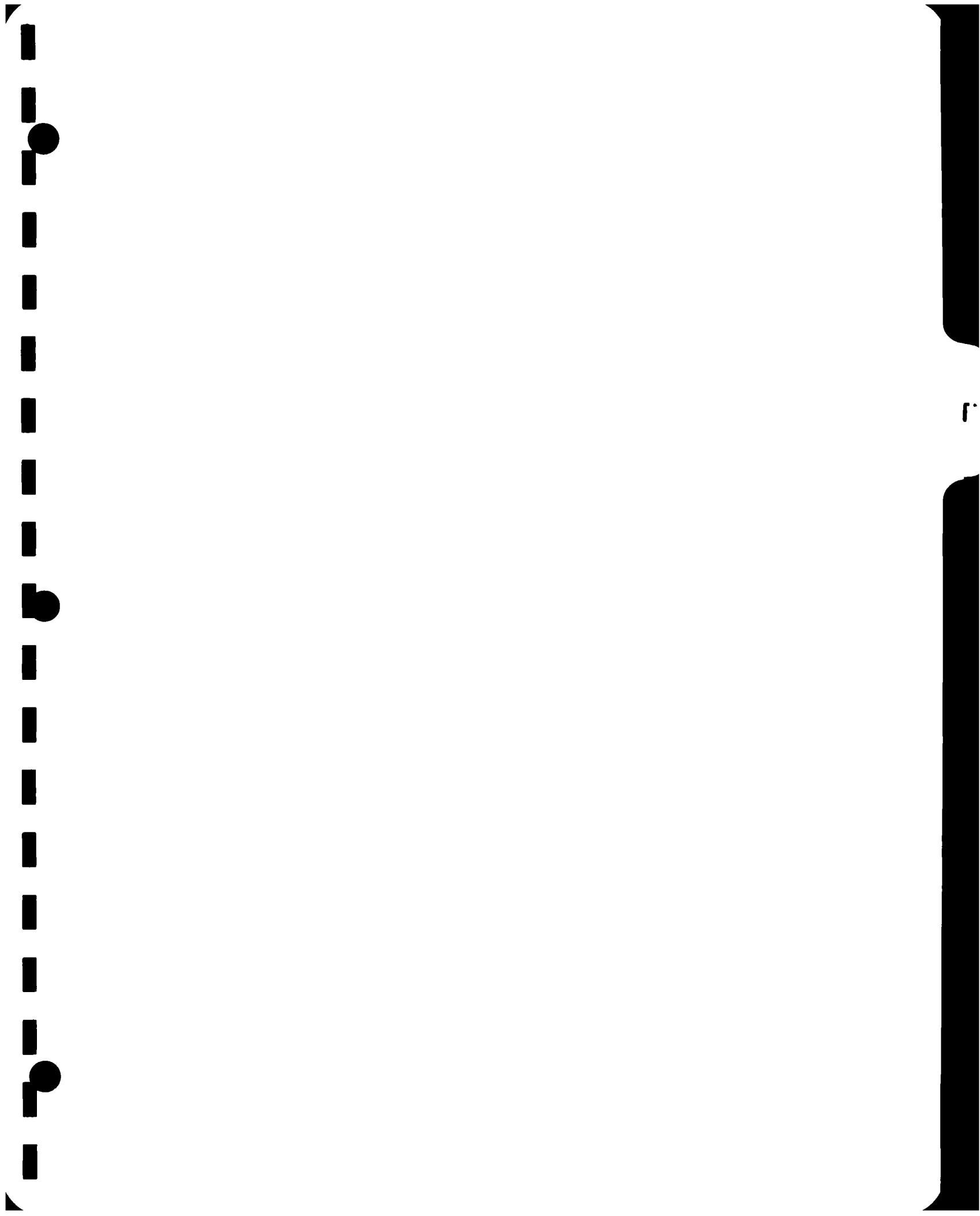
Blasting at the quarry may have created sufficient subsurface disturbance to cloud area ground water. Increased dissolved oxygen could be attributable to aeration and facilitate the growth of iron bacteria, as well.

Although no bacteria cultures were taken from the wells, the MPCA staff feels iron bacteria is the most probable cause of the poor quality of ground water at the Hader Ground Water Contamination Site, as opposed to a contaminant source.

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A



SITE NAME : HADER 64

SAMPLE LOCATION : 4-3

CRL - 287116507

SAMPLER : D LOFGREN

DATE 5/25/81 TIME 1405

**SITE NAME :** HADER SWCA

**SAMPLE LOCATION :** H-5

**CRL # :** E8YLUG506

**SAMPLER :** B LIFgren

**DATE :** 3/30/88 **TIME :** 1500

BB 3 30

SITE NAME :

SAMPLE LOCATION :

CRL :

SAMPLER :

DATE :            TIME : 59

88-3-10

SITE NAME : HADER GID  
SAMPLE LOCATION : H-7  
CRL : 881 DDOS  
SAMPLER : B. LOFLORPEN  
DATE : 3/3 / TIME : 1200

E  
SITE NAME HADER  
SAMPLE LOCATION S-2  
CRL - 88YLLD02  
SAMPLER B. L. GATE  
DATE 3/5/88 TIME 1200

88

SITE NAME: HAZZ  
SAMPLE LOCATION: H.V.  
CRL - 191-255  
SAMPLER: B. LIFSON  
DATE: 3-5-87 TIME:

ITEM NAME: ~~WALL~~  
SAMPLE LOCATION: ~~WALL~~  
CBL - ~~80112000~~  
NUMBER: ~~B-1000A~~  
DATE: ~~3/10/00~~

SITE NAME:   
SAMPLE LOCATION:   
CBL P#   
SAMPLER:   
DATE: TIME:

88 3 29

DATE : TIME :



SITE NAME : INTER GW CNT

SAMPLE LOCATION : H3

CRL # : 116502

SAMPLER : L - REN

DATE : 88-3-30 TIME : 1030

88 3 30

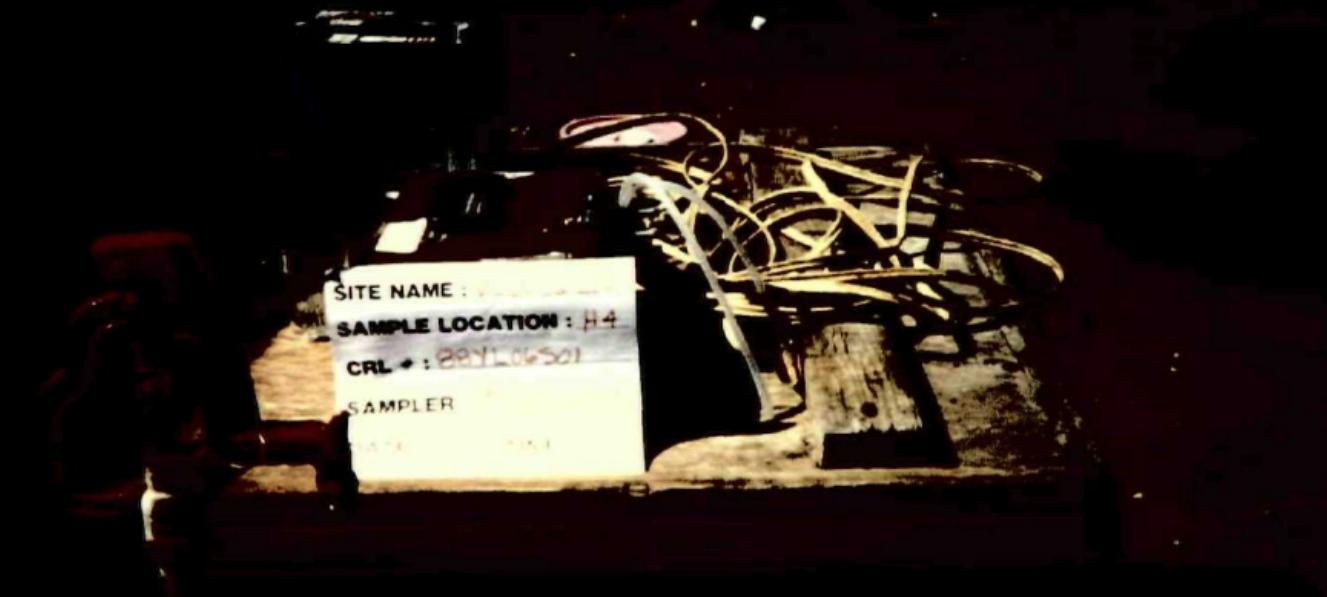


88 3 30

SITE NAME :  
SAMPLE LOCATION :  
CRL = :  
SAMPLER :  
DATE :            TIME :



- 88 3 3



88 3 30

SITE NAME HADER 2W

SAMPLE LOCATION H-2

TIME 10:00 AM

DATE 10/10/88

ANALYST J. D. COOPER

TEST TIME 10:00 AM

88 3 31



88 33:



Site Hader Groundwater Contamination

EPA # MND981961873

Date March 30, 1988

Time 12:56 a.m. p.m.

Direction Northwest

Weather Clear

Photographed by: Shawn Ruotsi

Sample ID # 88YL06S01

Description Well Head

Sample for S01



88 3 30

Site Hader Groundwater Contamination

EPA # MND981961873

Date March 30, 1988

Time 12:56 a.m. p.m.

Direction Northwest

Weather Clear

Photographed by: Shawn Ruotsi

Sample ID# 88YL06S01

Description Distance

Photo of S01



88 3 30

Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 10:30 a.m. p.m.

Direction Northwest

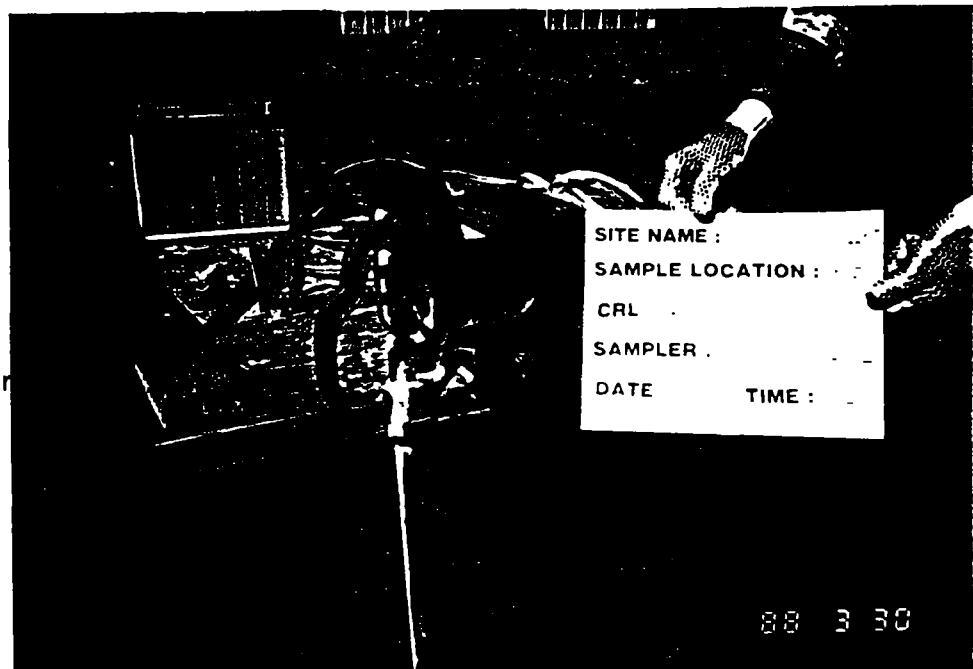
Weather Clear

Photographed by: Shawn Ruotsim

Sample ID # 88YL06S02

Description Wellhead

Sample of S02



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Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 10:30 a.m. p.m.

Direction Northwest

Weather Clear

Photographed by: Shawn Ruotsim

Sample ID# 88YL06S02

Description Distance Photo

of sample S02



Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 10:55 a.m. p.m.

Direction Northwest

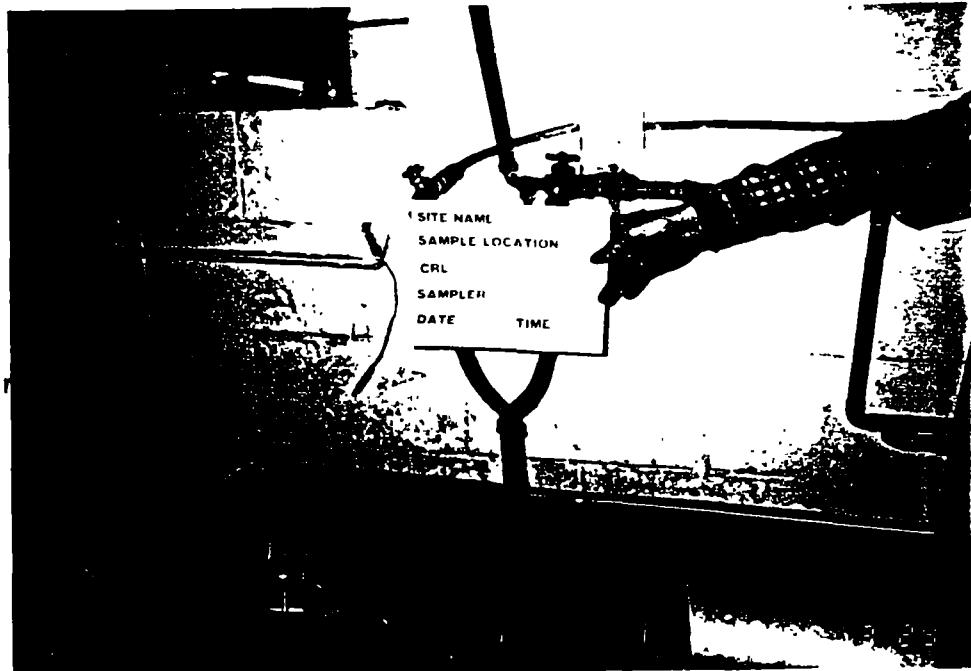
Weather Clear

Photographed by: Shawn Ruotsim

Sample ID # 88YL06S03

Description Sample taken

from faucet in milk barn



Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 10:55 a.m. p.m.

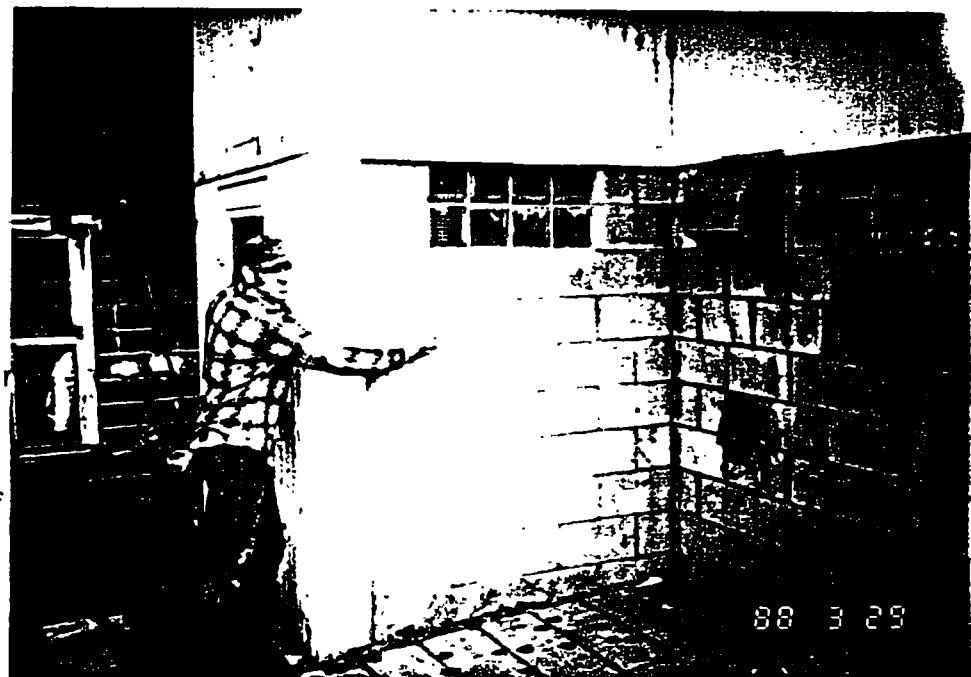
Direction Northwest

Weather Clear

Photographed by: Shawn Ruotsim

Sample ID# 88YL06S03

Description Distance photo of  
sample S03 at milk barn



Site Hader GW Contamination

EPA # MND981961873

Date March 29, 1988

Time 2:05 a.m. (p.m.)

Direction Northwest

Weather Clear

Photographed by: Shawn Ruotsi

Sample ID # 88YL06S04

Description Sample S04

taken at kitchen faucet



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Site Hader GW Contamination

EPA # MND981961873

Date March 29, 1988

Time 2:05 a.m. (p.m.)

Direction Northwest

Weather Clear

Photographed by: Shawn Ruotsi

Sample ID# 88YL06S04

Description Distance photo of  
sample S04 house



Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 10:52 a.m. p.m.

Direction Northwest

Weather Clear

Photographed by: Susan Price

Sample ID # 88YL06S05

Description Kitchen faucet

sample S05



BB 3 3 1

Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 10:52 a.m. p.m.

Direction Northwest

Weather Clear

Photographed by: Susan Price

Sample ID# 88YL06S05

Description Distance photo

of rear door at sample S05



BB 3 3 1

Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 12:06 a.m. p.m.

Direction Northwest

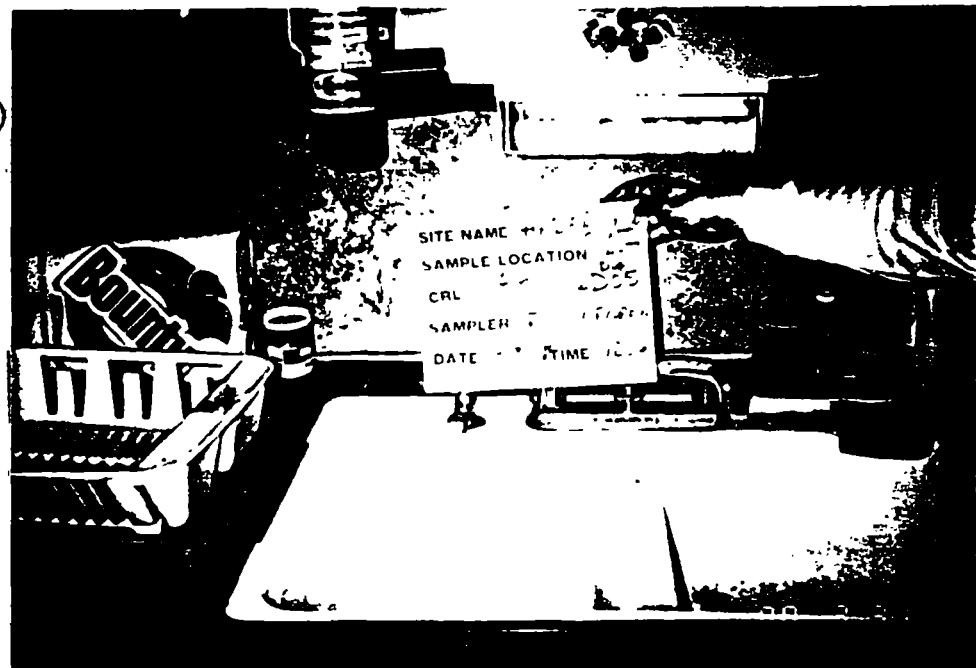
Weather Clear

Photographed by: Susan Price

Sample ID # 88YL06D05

Description Duplicate of S04

at kitchen sink, D05



Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 12:06 a.m. p.m.

Direction Northwest

Weather Clear

Photographed by: Susan Price

Sample ID# 88YL06D05

Description Distance Photo

of D04, duplicate of S04



Site Hader GW Contamination

EPA # MND9819661873

Date March 30, 1988

Time 3:00 a.m. p.m.

Direction Northwest

Weather Clear

Photographed by: Susan Price

Sample ID # 88YL06S06

Description Well head of

sample S06



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Site Hader GW Contamination

EPA # MND981961873

Date March 30, 1988

Time 3:00 a.m. p.m.

Direction Northwest

Weather Clear

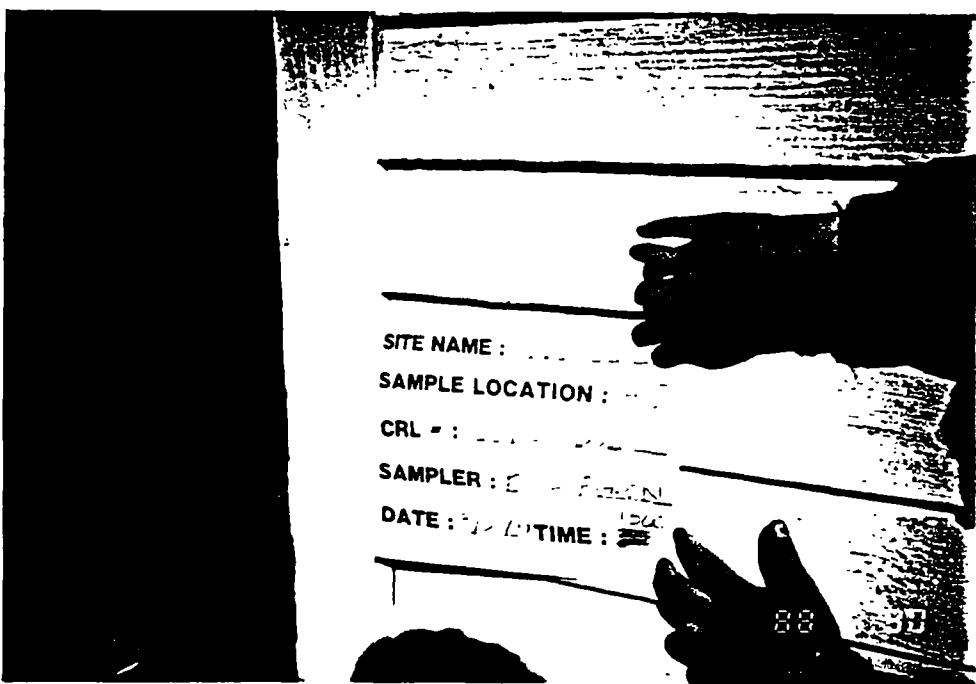
Photographed by: Susan Price

Sample ID# 88YL06S06

Description Distance photo

of sample S06 entrance of

well house



Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 2:05 a.m. p.m.

Direction North-Northwest

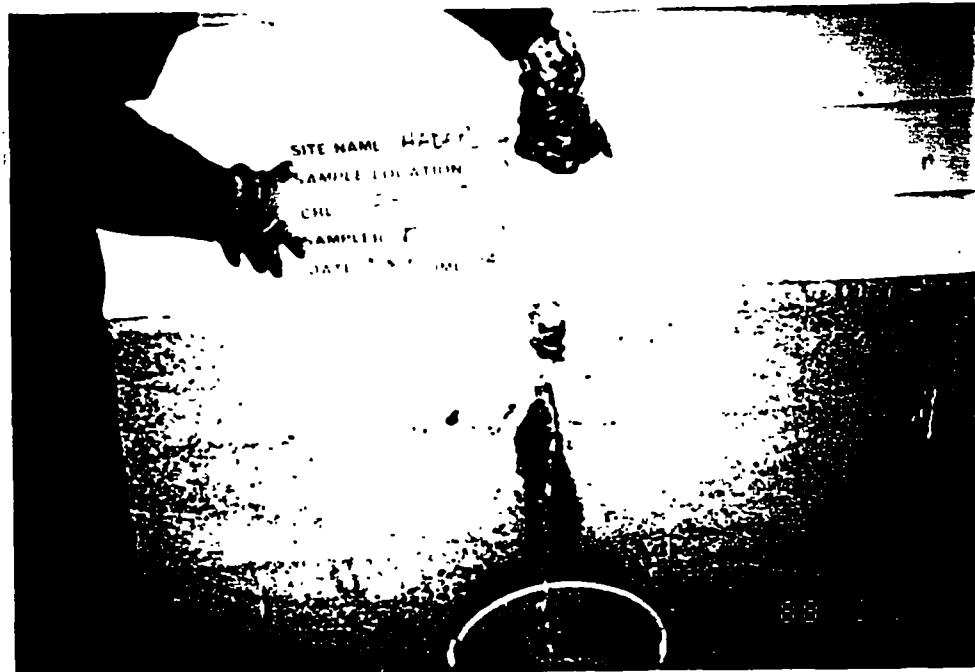
Weather Clear

Photographed by: Susan Price

Sample ID # 88YL06S07

Description Outdoor faucet

for sample S07



Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 2:05 a.m. p.m.

Direction North-Northwest

Weather Clear

Photographed by: Becky Lofgren

Sample ID# 88YL06S07

Description Distance photo

for sample S07



Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 3:05 a.m. p.m.

Direction North-Northwest

Weather Clear

Photographed by: Susan Price

Sample ID # 88YL06S08

Description Outdoor faucet

for sample S08



88 3 3 1

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Site Hader GW Contamination

EPA # MND981961873

Date March 31, 1988

Time 3:05 a.m. p.m.

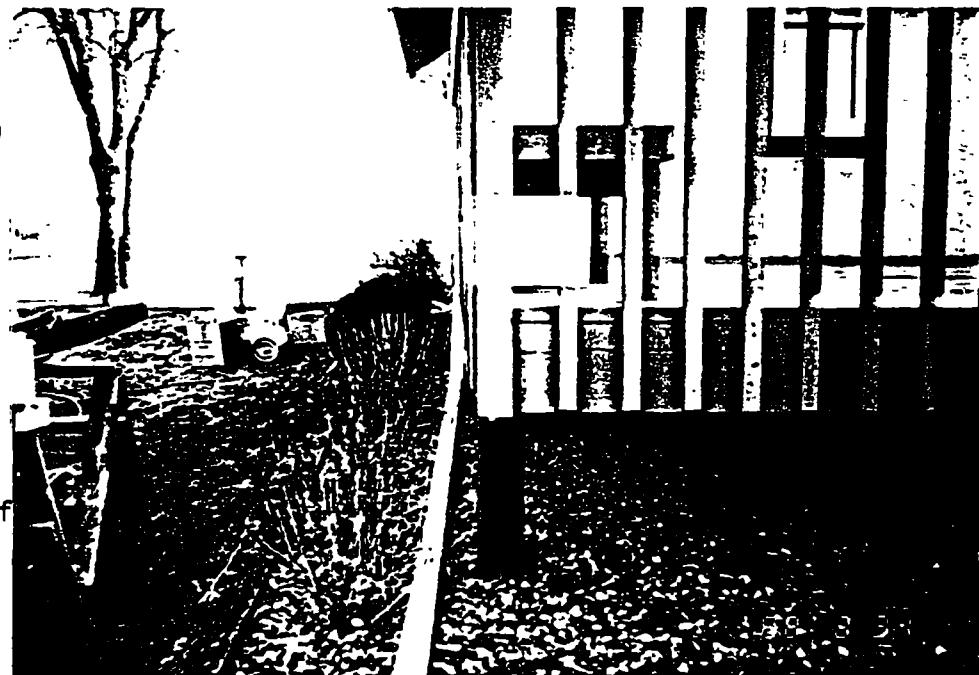
Direction North-Northwest

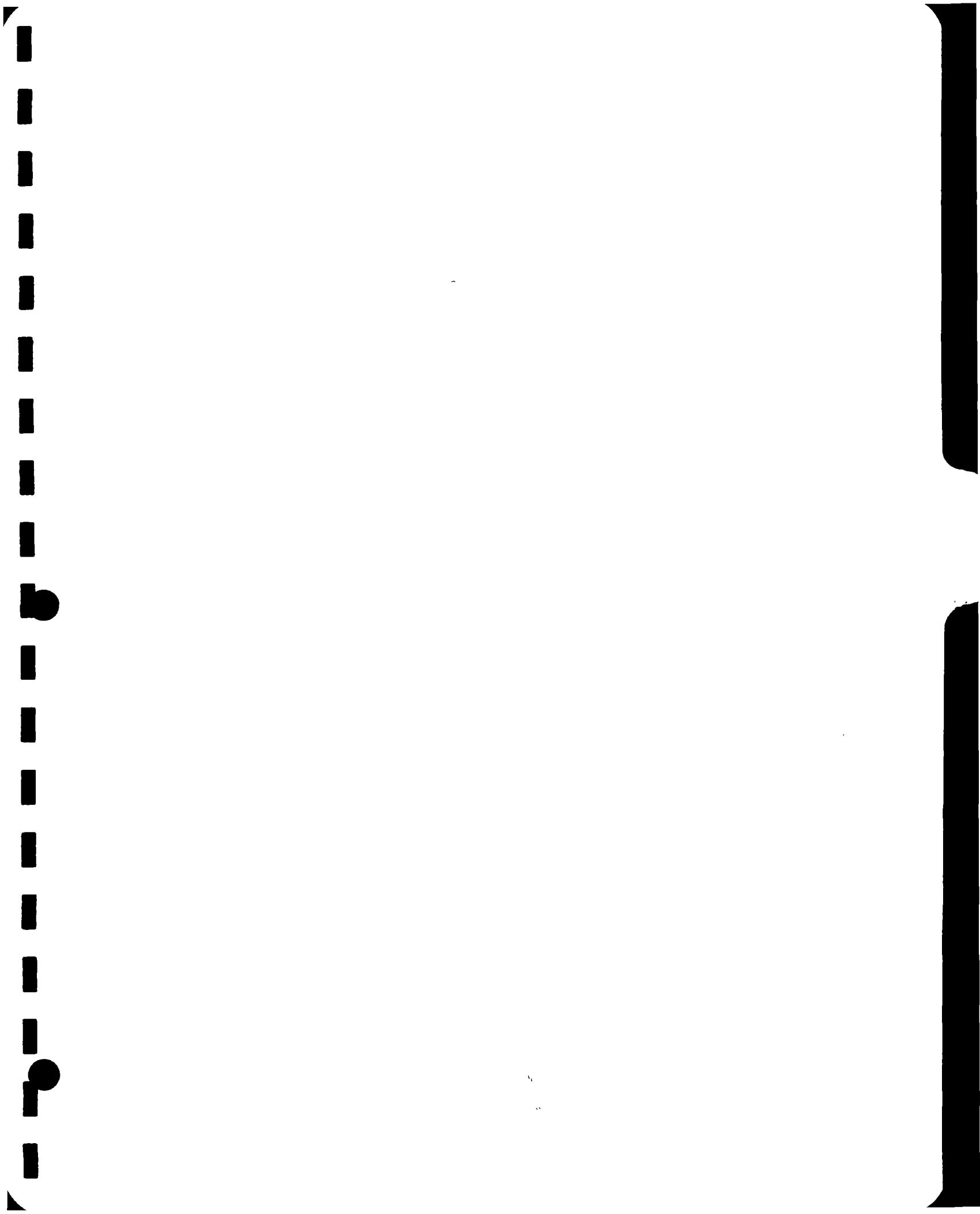
Weather Clear

Photographed by: Susan Price

Sample ID# 88YL06S08

Description Distance photo of  
sample S08







**Contract Laboratory Program**  
**Target Compound List**  
**Quantitation Limits**

COMPOUND	CAS #	WATER	SOIL SEDIMENT SLUDGE
Chloromethane	74-87-3	10 ug/L	10 ug/Kg
Bromomethane	74-83-9	10	10
Vinyl chloride	75-01-4	10	10
Chloroethane	75-00-3	10	10
Methylene chloride	75-09-2	5	5
Acetone	67-64-1	10	5
Carbon disulfide	75-15-0	5	5
1,1-dichloroethene	75-35-4	5	5
1,1-dichloroethane	75-34-3	5	5
1,2-dichloroethene (total)	540-59-0	5	5
Chloroform	67-66-3	5	5
1,2-dichloroethane	107-06-2	5	5
2-butanone (MEK)	78-93-3	10	10
1,1,1-trichloroethane	71-55-6	5	5
Carbon tetrachloride	56-23-5	5	5
Vinyl acetate	108-05-4	10	10
Bromodichloromethane	75-27-4	5	5
1,2-dichloropropane	78-87-5	5	5
cis-1,3-dichloropropene	10061-01-5	5	5
Trichloroethene	79-01-6	5	5
Dibromochloromethane	124-48-1	5	5
1,1,2-trichloroethane	79-00-5	5	5
Benzene	71-43-2	5	5
Trans-1,3-dichloropropene	10061-02-6	5	5
Bromoform	75-25-2	5	5
4-Methyl-2-pentanone	108-10-1	10	10
2-Hexanone	591-78-6	10	10
Tetrachloroethene	127-18-4	5	5
Tolene	108-88-3	5	5
1,1,2,2-tetrachloroethane	79-34-5	5	5
Chlorobenzene	108-90-7	5	5
Ethyl benzene	100-41-4	5	5
Styrene	100-42-5	5	5
Xylenes (total)	1330-20-7	5	5

**Contract Laboratory Program**  
**Target Compound List**  
**Semivolatiles Quantitation Limits**

COMPOUND	CAS #	WATER	SOIL SEDIMENT SLUDGE
Phenol	108-95-2	10 ug/L	330 ug/Kg
bis(2-Chloroethyl) ether	111-44-4	10	330
2-Chlorophenol	95-57-8	10	330
1,3-Dichlorobenzene	541-73-1	10	330
1,4-Dichlorobenzene	106-46-7	10	330
Benzyl Alcohol	100-51-6	10	330
1,2-Dichlorobenzene	95-50-1	10	330
2-Methylphenol	95-48-7	10	330
bis(2-Chloroisopropyl) ether	108-60-1	10	330
4-Methylphenol	106-44-5	10	330
N-Nitroso-di-n-dipropylamine	621-64-7	10	330
Hexachloroethane	67-72-1	10	330
Nitrobenzene	98-95-3	10	330
Isophorone	78-59-1	10	330
2-Nitrophenol	88-75-5	10	330
2,4-Dimethylphenol	105-67-9	10	330
Benzoic Acid	65-85-0	50	1600
bis(2-Chloroethoxy) methane	111-91-1	10	330
2,4-Dichlorophenol	120-83-2	10	330
1,2,4-Trichlorobenzene	120-82-1	10	330
Naphthalene	91-20-3	10	330
4-Chloroaniline	106-47-8	10	330
Hexachlorobutadiene	87-68-3	10	300
4-Chloro-3-methylphenol	59-50-7	10	330
2-Methylnaphthalene	91-57-6	10	330
Hexachlorocyclopentadiene	77-47-4	10	330
2,4,6-Trichlorophenol	88-06-2	10	330
2,4,5-Trichlorophenol	95-95-4	50	1600
2-Chloronaphthalene	91-58-7	10	330
2-Nitroaniline	88-74-4	50	1600
Dimethylphthalate	131-11-3	10	330
Acenaphthylene	208-96-8	10	330
2,6-Dinitrotoluene	606-20-2	10	330
3-Nitroaniline	99-09-2	50	1600
Acenaphthene	83-32-9	10	330
2,4-Dinitrophenol	51-28-5	50	1600
4-Nitrophenol	100-02-7	50	1600
Dibenzofuran	132-64-9	10	330
2,4-Dinitrotoluene	121-14-2	10	330
Diethylphthalate	84-66-2	10	330
4-Chlorophenyl-phenyl ether	7005-72-3	10	330

**Contract Laboratory Program**  
**Target Compound List**  
**Semivolatiles Quantitation Limits**

COMPOUND	CAS #	WATER	SOIL SLUDGE SEDIMENT
Fluorene	86-73-7	10 ug/L	330 ug/Kg
4-Nitroaniline	100-01-6	50	1600
4,6-Dinitro-2-methylphenol	534-52-1	50	1600
N-nitrosodiphenylamine	86-30-6	10	330
4-Bromophenyl-phenylether	101-55-3	10	330
Hexachlorobenzene	118-74-1	10	330
Pentachlorophenol	87-86-5	50	1600
Phenanthrene	85-01-8	10	330
Anthracene	120-12-7	10	330
Di-n-butylphthalate	84-74-2	10	330
Fluoranthene	206-44-0	10	330
Pyrene	129-00-0	10	330
Butylbenzylphthalate	85-68-7	10	330
3,3'-Dichlorobenzidine	91-94-1	20	660
Benzo(a)anthracene	56-55-3	10	330
Chrysene	218-01-9	10	330
bis(2-Ethylhexyl)phthalate	117-81-7	10	330
Di-n-octylphthalate	117-84-0	10	330
Benzo(b)fluoranthene	205-99-2	10	330
Benzo(k)fluoranthene	207-08-9	10	330
Benzo(a)pyrene	50-32-8	10	330
Indeno(1,2,3-cd)pyrene	193-39-5	10	330
Dibenz(a,h)anthracene	53-70-3	10	330
Benzo(g,h,i)perylene	191-24-2	10	330

**Contract Laboratory Program**  
**Target Compound List**  
**Pesticide and PCB Quantitation Limits**

COMPOUND	CAS #	WATER	SOIL SEDIMENT SLUDGE
alpha-BHC	319-84-6	0.05 ug/L	8 ug/Kg
beta-BHC	319-85-7	0.05	8
delta-BHC	319-86-8	0.05	8
gamma-BHC (Lindane)	58-89-9	0.05	8
Heptachlor	76-44-8	0.05	8
Aldrin	309-00-2	0.05	8
Heptachlor epoxide	1024-57-3	0.05	8
Endosulfan I	959-98-8	0.05	8
Dieldrin	60-57-1	0.10	16
4,4'-DDE	72-55-9	0.10	16
Endrin	72-20-8	0.10	16
Endosulfan II	33213-65-9	0.10	16
4,4'-DDD	72-54-8	0.10	16
Endosulfan sulfate	1031-07-8	0.10	16
4,4'-DDT	50-29-3	0.10	16
Methoxychlor (Mariate)	72-43-5	0.5	80
Endrin ketone	53494-70-5	0.10	16
alpha-Chlordane	5103-71-9	0.5	80
gamma-chlordane	5103-74-2	0.5	80
Toxaphene	8001-35-2	1.0	160
AROCLOR-1016	12674-11-2	0.5	80
AROCLOR-1221	11104-28-2	0.5	80
AROCLOR-1232	11141-16-5	0.5	80
AROCLOR-1242	53469-21-9	0.5	80
AROCLOR-1248	12672-29-6	0.5	80
AROCLOR-1254	11097-69-1	1.0	160
AROCLOR-1260	11096-82-5	1.0	160

**Contract Laboratory Program**  
**Target Analyte List**  
**Inorganic Quantitation Limits**

COMPOUND	PROCEDURE	SOIL WATER	SEDIMENT SLUDGE
Aluminum	ICP	200 ug/L	40 mg/Kg
Antimony	Furnace	60	2.4
Arsenic	Furnace	10	2
Barium	ICP	200	40
Beryllium	ICP	5	1
Cadmium	ICP	5	1
Calcium	ICP	5000	1000
Chromium	ICP	10	2
Cobalt	ICP	50	10
Copper	ICP	25	5
Iron	ICP	100	20
Lead	Furnace	5	1
Magnesium	ICP	5000	1000
Manganese	ICP	15	3
Mercury	Cold Vapor	0.2	0.008
Nickel	ICP	40	8
Potassium	ICP	5000	1000
Selenium	Furnace	5	1
Silver	ICP	10	2
Sodium	ICP	5000	1000
Thallium	Furnace	10	2
Vanadium	ICP	50	10
Zinc	ICP	20	4
Cyanide	Color	10	2

SPECIAL ANALYTICAL SERVICES DRINKING WATER  
VOLATILE QUANTITATION LIMITS

PARAMETER	CAS #	DETECTION LIMIT IN REAGENT WATER
Benzene	71-43-2	1.5 ug/L
Bromodichloromethane	74-27-4	1.5
Bromoform	75-25-2	1.5
Bromomethane	74-83-9	10
Carbon tetrachloride	56-23-5	1.5
Chlorobenzene	108-90-7	1.5
Chloroethane	75-00-3	1.5
2-Chloroethyl vinyl ether	110-75-8	1.5
Chloroform	67-66-3	1.5
Chloromethane	74-87-3	10
Dibromochloromethane	124-48-1	1.5
1,1-Dichloroethane	75-34-3	1.5
1,2-Dichloroethane	107-06-2	1.5
1,1-Dichloroethene	75-35-4	1.5
trans-1,2-Dichloroethene	156-60-5	1.5
1,2-Dichloropropane	78-87-5	1.5
cis-1,3-Dichloropropene	10061-01-5	2
trans-1,3-Dichloropropene	10061-02-6	1
Ethyl benzene	100-41-4	1.5
Methylene chloride *	75-09-2	1
1,1,2,2-Tetrachloroethane	79-34-5	1.5
Tetrachloroethene	127-18-4	1.5
Toluene *	108-88-3	1.5
1,1,1-Trichloroethane	71-55-6	1.5
1,1,2-Trichloroethane	79-00-5	1.5
Trichloroethene	79-01-6	1.5
Vinyl chloride	75-01-4	10
Acrolein	107-02-8	100
Acetone *	67-64-1	75
Acrylonitrile	107-13-1	50
Carbon disulfide	75-15-0	3
2-Butanone	78-93-3	(50)
Vinyl acetate	108-05-4	15
4-Methyl-2-pentanone	108-10-1	(3)
2-Hexanone	519-78-6	(50)
Styrene	100-42-5	1
m-Xylene	108-38-3	2
o-Xylene **	95-47-6	
p-Xylene **	106-42-3	2.5 **
Xylene (total)	1330-02-7	

\* Common laboratory solvents.

Blank limit is 5x method detection limit.

( ) Values in parentheses are estimates.

actual values are being determined at this time.

\*\* The o-xylene and p-xylene are reported as a total of the two.

SAS DRINKING WATER  
SEMIVOLATILES QUANTITATION LIMITS

PARAMETER	CAS #	DETECTION LIMIT
Aniline	62-53-3	1.5 ug/l
Bis(2-chloroethyl)ether	111-44-4	1.5
Phenol	108-95-2	2
2-Chlorophenol	95-57-8	2
1,3-Dichlorobenzene	541-73-1	2
1,4-Dichlorobenzene	106-46-7	2
1,2-Dichlorobenzene	95-50-1	2.5
Benzyl alcohol	100-51-6	2
Bis(2-chloroisopropyl)ether	39638-32-9	2.5
2-Methylphenol	95-48-7	1
Hexachloroethane	-67-72-1	2
n-Nitrosodipropylamine	621-64-7	1.5
Nitrobenzene	98-95-3	2.5
4-Methylphenol	88-75-5	1
Isophorone	78-59-1	2.5
2-Nitrophenol	88-75-5	2
2,4-Dimethylphenol	105-67-9	2
Bis(2-Chloroethoxy)methane	111-91-1	2.5
2,4-Dichlorophenol	120-83-2	2
1,2,4-Trichlorobenzene	120-82-1	2
Naphthalene	91-20-3	2
4-Chloroaniline	106-47-8	2
Hexachlorobutadiene	87-68-3	2.5
Benzoic Acid	65-85-0	(30)
2-Methylnaphthalene	91-57-6	2
4-Chloro-3-methylphenol	59-50-7	1.5
Hexachlorocyclopentadiene	77-47-4	2
2,4,6-Trichlorophenol	88-06-2	1.5
2,4,5-Trichlorophenol	95-95-4	1.5
2-Chloronaphthalene	91-58-7	1.5
Acenaphthylene	208-96-8	1.5
Dimethyl phthalate	131-11-3	1.5
2,6-Dinitrotoluene	606-20-2	1
Acenaphthene	83-32-9	1.5
3-Nitroaniline	99-09-2	2.5
Dibenzofuran	132-64-9	1
2,4-Dinitrophenol	51-28-5	(15)
2,4-Dinitrotoluene	121-14-2	1

SAS DRINKING WATER  
SEMIVOLATILE QUANTITATION LIMITS

PARAMETER	CAS #	DETECTION LIMIT
Fluorene	86-73-7	1 ug/L
4-Nitrophenol	100-02-7	1.5
4-Chlorophenyl phenyl ether	7005-72-3	1
Diethyl phthalate	84-66-2	1
4,6-Dinitro-2-methylphenol	534-52-1	(15)
1,2-Diphenylhydrazine	122-66-7	1
n-Nitrosodiphenylamine *	86-30-6	
Diphenylamine *	122-39-4	1.5
4-Nitroaniline	100-01-6	3
4-Bromophenyl-phenylether	101-55-3	1.5
Hexachlorobenzene	118-74-1	1.5
Pentachlorophenol	87-86-5	2
Phenanthrene	85-01-8	1
Anthracene	120-12-7	2.5
di-n-Butyl phthalate	84-74-2	2
Fluoranthene	206-44-0	1.5
Pyrene	129-00-0	1.5
Butyl benzyl phthalate	85-68-7	3.5
Chrysene **	218-01-9	
Benzo(A)Anthracene **	56-55-3	1.5
bis(2-ethylhexyl)phthalate	117-81-7	1
di-n-Octyl phthalate	117-84-0	1.5
Benzo(b)fluoranthene ***	205-99-2	
Benzo(k)fluoranthene ***	207-08-9	1.5
Benzo(a)pyrene	50-32-8	2
Indeno(1,2,3-cd)pyrene	193-39-5	3.5
Dibenzo(a,h)anthracene	53-70-3	2.5
Benzo(g,h,i)perylene	191-24-2	4
2-Nitroaniline .	88-74-4	1

\* These two parameters are reported as a total.

\*\* These two parameters are reported as a total.

\*\*\* These two parameters are reported as a total.

( ) Values in parentheses are estimates.

The actual values are being determined at this time.

Note: Limits are for reagent water.

**SAS DRINKING WATER  
PESTICIDE AND PCB QUANTITATION LIMITS**

<b>PARAMETER</b>	<b>CAS #</b>	<b>DETECTION LIMIT</b>
Aldrin	309-00-2	0.005 ug/L
alpha BHC	319-84-6	(0.010)
beta BHC	319-85-7	(0.005)
delta BHC	319-86-8	(0.005)
gamma BHC (Lindane)	58-89-9	0.005
Chlordane	57-74-9	(0.020)
4,4'-DDD	72-54-8	(0.020)
4,4'-DDE	72-55-9	(0.005)
4,4'-DDT	50-29-3	0.020
Dieldrin	60-57-1	0.010
Endosulfan I	959-98-8	0.010
Endosulfan II	33213-65-9	0.010
Endosulfan sulfate	1031-07-8	(0.10)
Endrin	72-20-8	0.010
Endrin Aldehyde	7421-93-4	(0.030)
Endrin Ketone	53494-70-5	(0.030)
Heptachlor	76-44-8	0.030
Heptachlor Epoxide	1024-57-3	0.005
4,4'-Methoxychlor	72-43-5	0.020
Toxaphene	8001-35-2	(0.25)
PCB-1242	53469-21-9	(0.10)
PCB-1248	12672-29-6	(0.10)
PCB-1254	11097-69-1	(0.10)
PCB-1260	11096-82-5	(0.10)

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( ) Values in parentheses are estimates.  
Actual values are being determined at this time.

Note: Limits are for reagent water.

SAS DRINKING WATER  
INORGANIC DETECTION LIMITS

JANUARY 1986

PARAMETER	PROCEDURE	DETECTION LIMIT
Aluminum	ICP	100
Antimony	GFAA	2
Arsenic	GFAA	2
Barium	ICP	50
Beryllium	ICP	5
Cadmium	ICP	10
Cadmium	GFAA	0.2
Calcium	ICP	1000
Chromium	ICP	10
Cobalt	ICP	10
Copper	ICP	10
Iron	ICP	100
Lead	GFAA	2
Magnesium	ICP	1000
Manganese	ICP	10
Mercury	Cold Vapor	0.2
Nickel	ICP	20
Potassium	ICP	2000
Selenium	GFAA	2
Silver	ICP	5
Sodium	ICP	1000
Thallium	GFAA	2
Tin	ICP	40
Vanadium	ICP	10
Zinc	ICP	20
Cyanide	Colorimetric	5.0

Note: The above list may or may not contain compounds that are routinely analyzed at CRL for low level detection limits for drinking water.

See inorganic Routine Analytical Services (RAS) for related CAS #.

If gel permeation chromatography, "GPC Cleanup" was performed, enter "Y" for yes. Otherwise, enter "N" for no, if GPC was not performed.

Enter pH for semivolatile and pesticides/PCBs, reported to 0.1 pH units.

"Date Received" is the date of sample receipt at the laboratory, as noted on the Traffic Report (i.e., the VTSR). It should be entered as MM/DD/YY.

"Date Extracted" and "Date Analyzed" should be entered in a similar fashion. For pesticide/PCB samples, the date of analysis should be the date of the first GC analysis performed. The date of sample receipt will be compared with the extraction and analysis dates of each fraction to ensure that contract holding times were not exceeded.

If a sample has been diluted for analysis, enter the "Dilution Factor" as a decimal number, such as 0.001 for a 1 to 1000 dilution of the sample. If a sample was not diluted, enter 1.

For positively identified TCL compounds, the contractor shall report the concentrations detected as uncorrected for blank contaminants.

For volatile and semivolatile results, report analytical results to one significant figure if the value is less than 10, and two significant figures above 10.

Report all pesticides/PCB results to two significant figures.

The appropriate concentration units, ug/L or ug/kg, must be entered.

If the result is a value greater than or equal to the quantitation limit, report the value.

Under the column labeled "Q" for qualifier, flag each result with the specific Data Reporting Qualifiers listed below. The Contractor is encouraged to use additional flags or footnotes. The definition of such flags must be explicit and must be included in the Case Narrative.

For reporting results to the USEPA, the following contract specific qualifiers are to be used. The seven qualifiers defined below are not subject to modification by the laboratory. Up to five qualifiers may be reported on Form I for each compound.

The seven EPA-defined qualifiers to be used are as follows:

U - Indicates compound was analyzed for but not detected. The sample quantitation limit must be corrected for dilution and for percent moisture. For example, 10 U for phenol in water if the sample final volume is the protocol-specified final volume. If a 1 to 10 dilution of extract is necessary, the reported limit is 100 U. For a soil sample, the value must also be adjusted for percent moisture. For example, if the sample had 24% moisture

and a 1 to 10 dilution factor, the sample quantitation limit for phenol (330 U) would be corrected to:

$$\frac{(330 \text{ U}) \times df}{D} \quad \text{where } D = \frac{100 - \% \text{ moisture}}{100}$$

and df = dilution factor

$$\text{at 24\% moisture, } D = \frac{100-24}{100} = 0.76$$

$$\frac{(330 \text{ U}) \times 10}{.76} = 4300 \text{ U rounded to the appropriate number of significant figures}$$

For soil samples subjected to GPC clean-up procedures, the CRQL is also multiplied by 2, to account for the fact that only half of the extract is recovered.

- J - Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed, or when the mass spectral data indicate the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than zero. For example, if the sample quantitation limit is 10 ug/L, but a concentration of 3 ug/L is calculated, report it as 3J. The sample quantitation limit must be adjusted for both dilution and percent moisture as discussed for the U flag, so that if a sample with 24% moisture and a 1 to 10 dilution factor has a calculated concentration of 300 ug/L and a sample quantitation limit of 430 ug/kg, report the concentration as 300J on Form I.
- C - This flag applies to pesticide results where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10 \text{ ng}/\mu\text{l}$  in the final extract shall be confirmed by GC/MS.
- B - This flag is used when the analyte is found in the associated blank as well as in the sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action. This flag must be used for a TIC as well as for a positively identified TCL compound.
- E - This flag identifies compounds whose concentrations exceed the calibration range of the GC/MS instrument for that specific analysis. This flag will not apply to pesticides/PCBs analyzed by GC/EC methods. If one or more compounds have a response greater than full scale, the sample or extract must be diluted and re-analyzed according to the specifications in Exhibit D. All such compounds with a response greater than full scale should have the concentration flagged with an "E" on the Form I for the original analysis. If the dilution of the extract causes any compounds identified in the first analysis to be below the calibration range in the second analysis, then the results of both analyses shall be reported on separate Forms I. The Form I for the diluted sample shall have the "DL" suffix appended to the sample number.

- D - This flag identifies all compounds identified in an analysis at a secondary dilution factor. If a sample or extract is re-analyzed at a higher dilution factor, as in the "E" flag above, the "DL" suffix is appended to the sample number on the Form I for the diluted sample, and all concentration values reported on that Form I are flagged with the "D" flag.
- A - This flag indicates that a TIC is a suspected aldol-condensation product.
- X - Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the Sample Data Summary Package and the Case Narrative. If more than one is required, use "Y" and "Z", as needed. If more than five qualifiers are required for a sample result, use the "X" flag to combine several flags, as needed. For instance, the "X" flag might combine the "A", "B", and "D" flags for some sample.

The combination of flags "BU" or "UB" is expressly prohibited. Blank contaminants are flagged "B" only when they are also detected in the sample.

If analyses at two different dilution factors are required (see Exhibit D), follow the data reporting instructions given in Exhibit D and with the "D" and "E" flags above.

## 2. Form I VOA-TIC and Form I SV-TIC

Fill in all header information as above.

Report Tentatively Identified Compounds (TIC) including CAS number, compound name, retention time, and the estimated concentration (criteria for reporting TICs are given in Exhibit D, Section IV). Retention time must be reported in minutes and decimal minutes, not seconds or minutes:seconds.

If in the opinion of the mass spectral interpretation specialist, no valid tentative identification can be made, the compound shall be reported as unknown.

Include a Form I VOA-TIC or SV-TIC for every volatile and semivolatile fraction of every sample and method blank analyzed, even if no TICs are found. Total the number of TICs found, including aldol-condensation products (but see below), and enter this number in the "Number TICs found." If none were found, enter "0" (zero).

If the name of a compound exceeds the 28 spaces in the TIC column, truncate the name to 28 characters. If the compound is an unknown, restrict description to no more than 28 characters (i.e., unknown hydrocarbon, etc.).

Peaks that are suspected as aldol-condensation reaction products (i.e., 4-methyl-4-hydroxy-2-pentanone and 4-methyl-3-pentene-2-one) shall be summarized on this form, flagged "A", and included in the total "Number TICs found," but not counted as part of the 20 most intense non-TCL semi-volatile compounds to be searched.

C. Surrogate Recovery (Form II)

Form II is used to report the recoveries of the surrogate compounds added to each sample, blank, matrix spike, and matrix spike duplicate. Form II is matrix-specific as well as fraction-specific, so that surrogate recoveries for volatile water samples are reported on a different version of Form II than volatile soil sample surrogate recoveries.

Complete the header information and enter EPA Sample Numbers as described in part A. For soil samples only, specify the "level" as "LOW" or "MED", as on Form I. Do not mix low and medium level samples on one form. Complete one for each level. For each surrogate, report the percent recovery to the number of significant figures given by the QC limits at the bottom of the form.

Flag each surrogate recovery outside the QC limits with an asterisk (\*). The asterisk must be placed in the last space in each appropriate column, under the "#" symbol. In the far righthand column, total the number of surrogate recoveries outside the QC limits for each sample. If no surrogates were outside the limits, enter "0".

If the surrogates are diluted out in any analysis, enter the calculated recovery or "0" (zero) if the surrogate is not detected, and flag the surrogate recoveries with a "D" in the column under the "#" symbol. Do not include results flagged "D" in the total number of recoveries for each sample outside the QC limits.

The pesticide surrogate recovery limits are only advisory, but the contractor must flag those recoveries outside the advisory QC limits or diluted out, nonetheless.

Number all pages as described in part A.

D. Matrix Spike/Matrix Spike Duplicate Recovery (Form III)

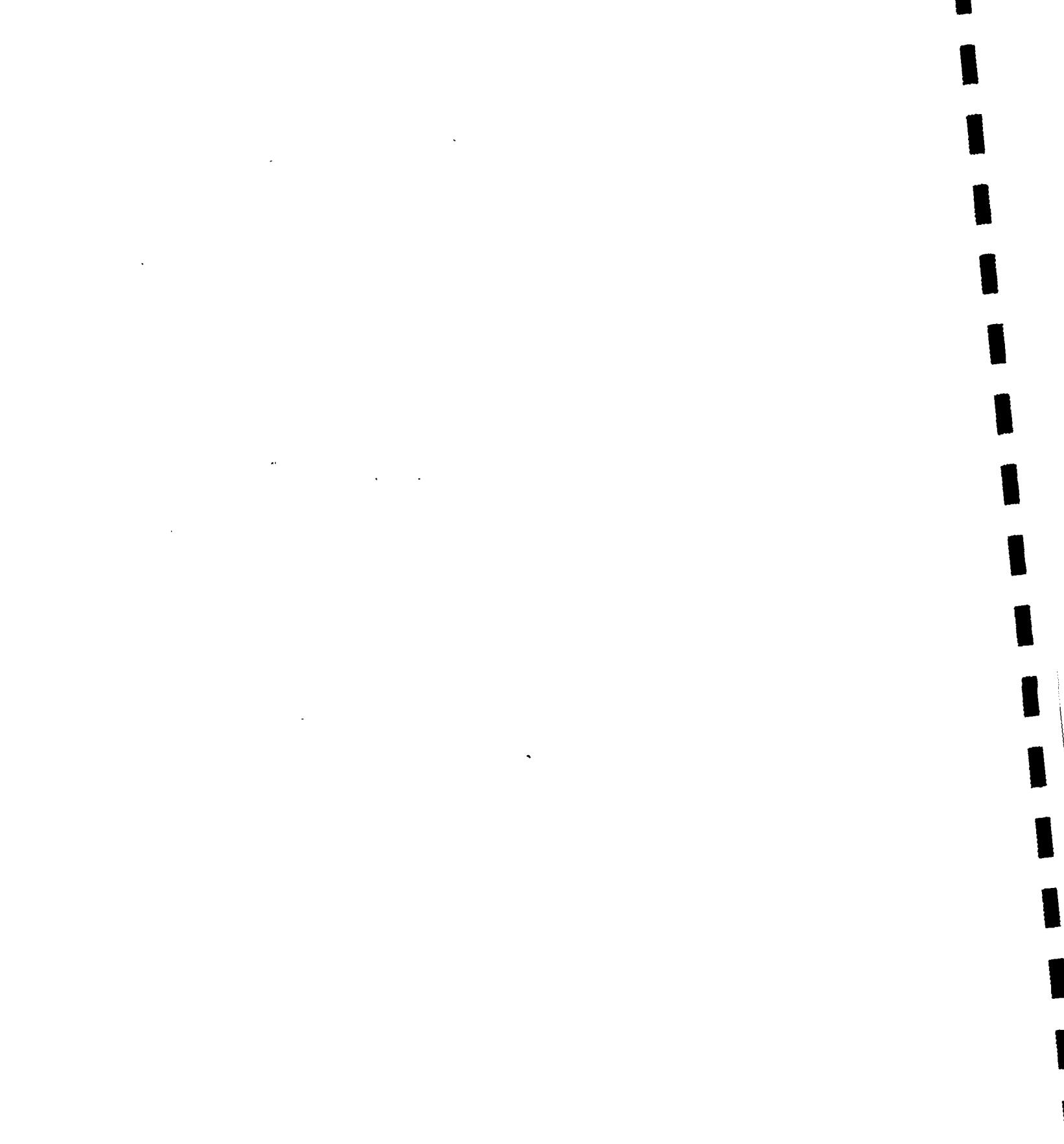
This form is used to report the results of the analyses of a matrix spike and matrix spike duplicate. As with the surrogate recovery form (II), the form is matrix-specific within each fraction.

Complete the header information as instructed in Part A, including the EPA Sample Number for the matrix spike without the suffixes MS or MSD.

For soil samples, specify "level" as "LOW" or "MED", as on Form I. Cases containing soil samples at both levels require MS/MSD at each level, therefore, for soils, prepare one form for each level.

All water samples are "Low". Therefore, there is no MS/MSD for "medium level waters", and none shall be reported.

# **ORGANIC DATA**



RECEIVED

MAY 12 88

MPCA, Ground Water  
& Solid Waste Div.

DATE: May 9, 1988

TO: Minnesota Pollution Control Agency  
Site Assessment Unit  
Program Development Section  
Groundwater and Solid Waste Division  
520 Lafayette Road  
St. Paul, MN 55155

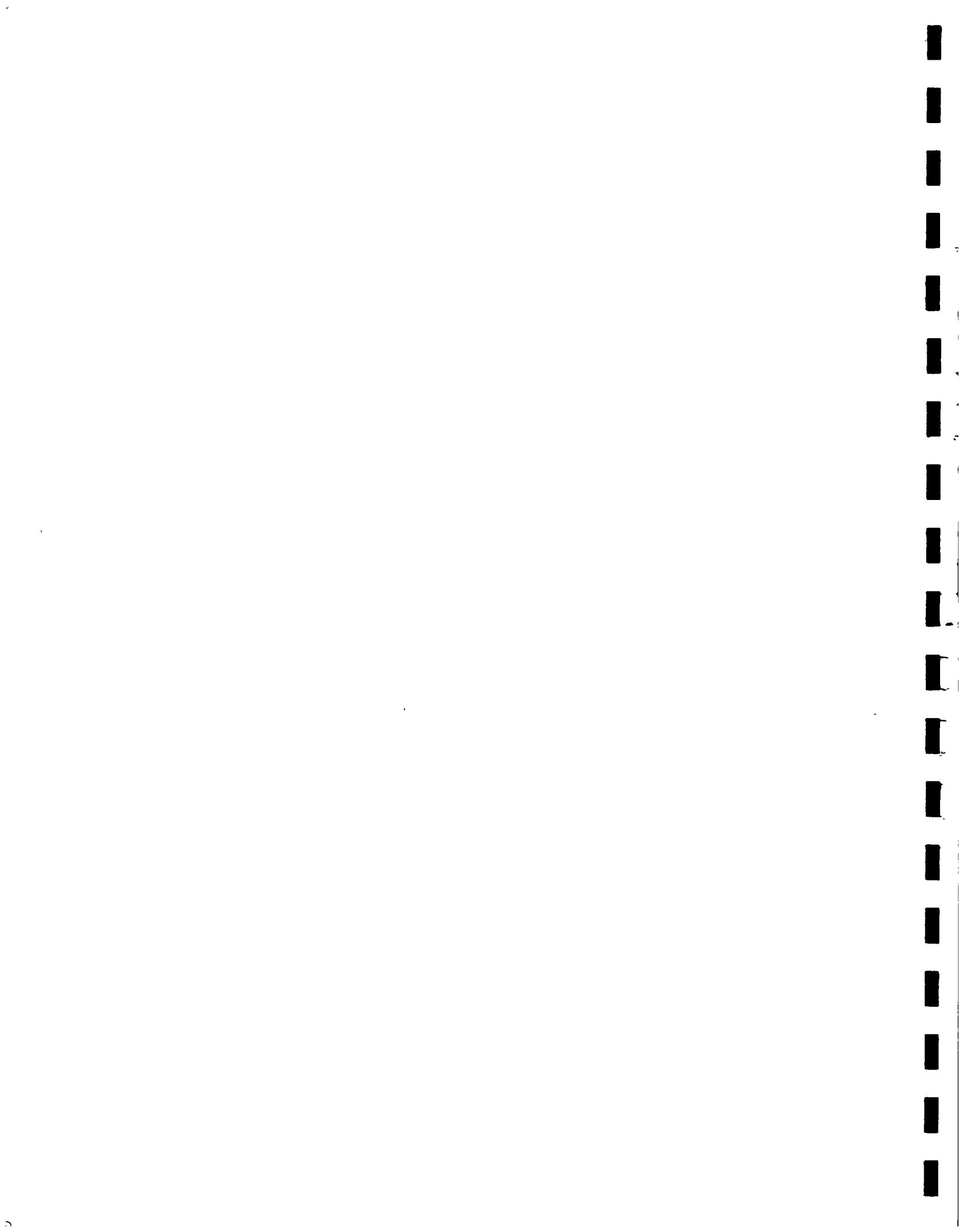
ATTN: Doug Day, Supervisor

Case No.	Site	Contract Lab	SF No.	No. Samples
9288SAS3554E	HADER GROUND WATER CONT.	S-CUBED	SF5011	10

Enclosed is one REVIEWED case. Please DO NOT return it.

FROM: U.S. EPA  
Region V  
Central Regional Laboratory  
556 S. Clark, 10th Floor  
Chicago, IL 60605

SENT BY: *Erlinda Luz M. Arreola*  
WESTON - ESAT



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

SITE: S-6-88

SUBJECT: Review of Region V CLP Data  
Received for Review on 4-25-88FROM: Curtis Ross, Director (5SCR) *Patricia J. Chumbley fm*  
Central Regional LaboratoryTO: Data User: MPCA  
DUANE KOVSKY, MPCA

We have reviewed the data for the following case(s).

SITE NAME: HADER GROUND WATER (MW) SMO case No. 9288SAS3554E  
EPA Data Set No. SF5011 No. of Samples: 10 D.U./Activity Numbers Y905/C72277CRL No. 88YLO6S01-S08, 005, R01SMO Traffic No. ET831-840CLP Laboratory: SCUBED Hrs. Required for Review: 7

Following are our findings:

This review covers the analysis of 10 low concentration water samples. These were analyzed for semi-volatiles, volatiles and pesticide/PCBs. One sample (ET840) was analyzed for volatiles only.

Several problems which should be noted by the data user for these samples are outlined on the following pages.

*John R. Stenger  
ESAT- WESTON  
5-8-88*

- Data are acceptable for use.
- Data are acceptable for use with qualifications noted above.
- Data are preliminary - pending verification by Contractor Laboratory.
- Data are unacceptable.

cc: Duane Geuder, Quality Assurance Officer, EPA Support Services  
James Petty, Chief Quality Assurance Research, EMSL, Las Vegas

## DATA QUALIFIERS

Contractor: S-CUBED

Case 9288 SAS 3554E

Below is a summary of the out-of-control audits and the possible effect on the data for this case:

1) Holding Time

The extractions for both the semi-volatile and pesticide/PCBs fractions for samples ET831 thru ET839 exceeded the 7 days from sampling date. Holding time - All results from these fractions should be considered estimated. (Positive (P), Negative (N)).

2) Tuning

The GC/MS tuning criteria was met for both the volatile and semi-volatile fractions. Overall instrument performance was acceptable. All instrument performance criteria were met for the pesticide/PCB fraction.

3) Calibration

Calibration outliers for the volatile and semi-volatile fractions are noted on the calibration outlier sheets. Calibration criteria for the pesticide/PCB fraction were acceptable.

4) Blanks

Common contaminants were found in the volatile and semi-volatile blanks. Non-HSL contaminants were found in the semi-volatile blank. No contaminants were found in the pesticide/PCB blank. All contaminants were at acceptable levels.

5) Surrogate Recoveries

Recoveries for Phenol-05 and 2-Fluorophenol in the semi-volatile fractions of samples ET834, ET835 and ET836 were less than 10% and under the QC limits. The fractions were re-injected with similar results. There was insufficient sample to re-extract.

Reviewed by:

John R. Marcyho

Phone:

✓ 312-353-2917

Date:

5-2-88

## DATA QUALIFIERS

Contractor: S-CUBEDCase 9288 SAS 3554E

Below is a summary of the out-of-control audits and the possible effect on the data for this case:

For these fractions of the three mentioned samples, positive results should be flagged estimated (I) and negative results flagged unusable (R).

6) Matrix Spike (MS)

The % recoveries for Endrin are above the QC limit for both the matrix spike (MS). Therefore for sample ET831 the results for Endrin should be flagged estimated (positive (I)), negative (N)).

7) Field Duplicate

Samples ET835 and ET836 were duplicates. This was relative agreement between these samples.

8) Compound Identification

All compound identifications were acceptable.

The problems identified in this review will not directly affect the quality and soundness of the data as a whole. Therefore, I find the data contained within acceptable for use.

Reviewed by:

John R. Haynes

Phone:

312-353-6917

## UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V

## CALIBRATION OUTLIERS

## VOLATILE HSL COMPOUNDS

CASE/SAS # 9288 SAS35SYECONTRACTOR S CUBED

Instrument # <u>VGA</u>	Init. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.
DATE/TIME:	3-28-88	4-4 0913	4-4 2016				
	RF %RSD *	RF %D *	RF %D *	RF %D *	RF %D *	RF %D *	RF %D *
Chloromethane							
Bromomethane							
Vinyl Chloride							
Chloroethane							
Methylene Chloride							
Acetone							
Carbon Disulfide							
1,1-Dichloroethane							
1,1-Dichloroethene							
Trans-1,2-Dichloroethene							
Chloroform							
2-Butanone	0.0187 J/R	0.01180 J/R	0.01170 J/R				
1,2-Dichloroethane							
1,1,1-Trichloroethane							
Carbon Tetrachloride							
Vinyl Acetate							
Bromodichloromethane							
1,2-Dichloropropane							
Trans-1,3-Dichloropropene							
Trichloroethene							
Dibromochloromethane							
1,1,2-Trichloroethane							
Benzene							
cis-1,3-Dichloropropene							
2-Chloroethylvinylether							
Bromoform							
4-Methyl-2-Pentanone							
2-Hexanone							
Tetrachloroethene							
1,1,2,2-Tetrachloroethane							
Toluene							
Chlorobenzene							
Ethylbenzene							
Styrene							
m-Xylene							
o/p-Xylene							
ACROLEIN	0.0393.33 J/R	0.049170 J/R	0.02723.0 J/R				
AFFECTED SAMPLES:		LBLK 44V.1021	LBLK 44V.1021				
Reviewer's Initials/Date: <u>JG 5-8-88</u>		ET831	ET837				
		ET833	ET831MS				
		ET832	ET831MSN				
		ET831	MATRIX BLANK				
		ET835					
		ET836					
		ET838					
		ET839					
		ET840					

\* These flags should be applied to the analytes on the sample data sheets.

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V  
 CALIBRATION OUTLIERS  
 SEMIVOLATILE HSL COMPOUNDS

(Page 1)

CASE/SAS # 9288 SAS 3554 ECONTRACTOR S-CUBED

Instrument # HP#1	Init. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.						
DATE/TIME:	4-10-88											
	RF	%RSD *	RF	%D *	RF	%D *	RF	%D *	RF	%D *	RF	%D *
Phenol												
bis(-2-Chloroethyl)Ether												
2-Chlorophenol												
1,3-Dichlorobenzene												
1,4-Dichlorobenzene												
Benzyl Alcohol												
1,2-Dichlorobenzene												
2-Methylphenol												
bis(2-chloroisopropyl)Ether												
4-Methylphenol												
N-Nitroso-Di-n-Propylamine												
Hexachloroethane												
Nitrobenzene												
Isophorone												
2-Nitrophenol												
2,4-Dimethylphenol												
Benzoic Acid												
bis(2-Chloroethoxy)Methane												
2,4-Dichlorophenol												
1,2,4-Trichlorobenzene												
Naphthalene												
4-Chloroaniline												
Hexachlorobutadiene												
4-Chloro-3-Methylphenol												
2-Methylnaphthalene												
Hexachlorocyclopentadiene												
2,4,6-Trichlorophenol												
2,4,5-Trichlorophenol												
2-Chloronaphthalene												
2-Nitroaniline												
Dimethyl Phthalate												
Acenaphthylene												
3-Nitroaniline												
Acenaphthene												
2,4-Dinitrophenol												
4-Nitrophenol												
Dibenzofuran												
AFFECTED SAMPLES:	LAB PLANK 47											
	ETP33											
	ETP34											
Reviewer												
Initials/Date:	5-2-88											

\* These flags should be applied to the analytes on the sample data sheets.

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V  
 CALIBRATION OUTLIERS  
 SEMIVOLATILE HSL COMPOUNDS

Page 2

CASE/SAS # 9288 SAS 3554 ECONTRACTOR S-CUBED

Instrument # HP# /	Init. Cal.	Cont. Cal.								
DATE/TIME:	4-10-88									
	RF	%RSD *	RF	%D *	RF	%D *	RF	%D *	RF	%D *
2,4-Dinitrotoluene										
2,6-Dinitrotoluene										
Diethylphthalate										
4-Chlorophenyl-phenylether										
Fluorene										
4-Nitroaniline										
4,6-Dinitro-2-Methylphenol										
N-Nitrosodiphenylamine										
4-Bromophenyl-phenylether										
Hexachlorobenzene										
Pentachlorophenol										
Phenanthrene										
Anthracene										
Di-n-Butylphthalate										
Fluoranthene										
Pyrene										
Butylbenzylphthalate										
Benzo(a)Anthracene										
bis(2-Ethylhexyl)Phthalate										
Chrysene										
Di-n-Octyl Phthalate										
Benzo(b)Fluoranthene										
Benzo(k)Fluoranthene										
Benzo(a)Pyrene										
Indeno(1,2,3-cd)Pyrene										
Dibenz(a,h)Anthracene										
Benzo(g,h,i) Perylene										

SEE PAGE 1 FOR AFFECTED SAMPLES.

\* These flags should be applied to the analytes on the sample data sheets.

Reviewer's Initials/Date:

*JB* 5-2-88

8/87

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V  
 CALIBRATION OUTLIERS  
 SEMIVOLATILE HSL COMPOUNDS

(Page 1)

CASE/SAS # 9288SAS3554ECONTRACTOR S-CUBED

Instrument # HP # 1	Init. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.	Cont. Cal.
DATE/TIME:	4-13-88					
Phenol						
bis(-2-Chloroethyl)Ether						
2-Chlorophenol						
1,3-Dichlorobenzene						
1,4-Dichlorobenzene						
Benzyl Alcohol						
1,2-Dichlorobenzene						
2-Methylphenol						
bis(2-chloroisopropyl)Ether						
4-Methylphenol						
N-Nitroso-Di-n-Propylamine						
Hexachloroethane						
Nitrobenzene						
Isophorone						
2-Nitrophenol						
2,4-Dimethylphenol						
Benzoic Acid						
bis(2-Chloroethoxy)Methane						
2,4-Dichlorophenol						
1,2,4-Trichlorobenzene						
Naphthalene						
4-Chloroaniline						
Hexachlorobutadiene						
4-Chloro-3-Methylphenol						
2-Methylnaphthalene						
Hexachlorocyclopentadiene						
2,4,6-Trichlorophenol						
2,4,5-Trichlorophenol						
2-Chloronaphthalene						
2-Nitroaniline						
Dimethyl Phthalate						
Acenaphthylene						
3-Nitroaniline						
Acenaphthene						
2,4-Dinitrophenol						
4-Nitrophenol						
Dibenzofuran						
AFFECTED SAMPLES:	RRPAG ET 837 ET 838					
Reviewer Initials/Date:	<u>JM</u> 5-9-88					

\* These flags should be applied to the analytes on the sample data sheets.

8/87

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V  
 CALIBRATION OUTLIERS  
 SEMIVOLATILE HSL COMPOUNDS

Page 2

CASE/SAS # 9288 SAS3554ECONTRACTOR S-CUGED

Instrument # HP #/	Init. Cal.	Cont. Cal.								
DATE/TIME:	4-13-88									
	RF	%RSD *	RF	%D *	RF	%D *	RF	%D *	RF	%D *
2,4-Dinitrotoluene										
2,6-Dinitrotoluene										
Diethylphthalate										
4-Chlorophenyl-phenylether										
Fluorene										
4-Nitroaniline										
4,6-Dinitro-2-Methylphenol										
N-Nitrosodiphenylamine										
4-Bromophenyl-phenylether										
Hexachlorobenzene										
Pentachlorophenol										
Phenanthrene										
Anthracene										
Di-n-Butylphthalate										
Fluoranthene										
Pyrene										
Butylbenzylphthalate										
Benzo(a)Anthracene										
bis(2-Ethylhexyl)Phthalate										
Chrysene										
Di-n-Octyl Phthalate										
Benzo(b)Fluoranthene										
Benzo(k)Fluoranthene										
Benzo(a)Pyrene										
Indeno(1,2,3-cd)Pyrene										
Dibenz(a,h)Anthracene										
Benzo(g,h,i) Perylene										

SEE PAGE 1 FOR AFFECTED SAMPLES.

\* These flags should be applied to the analytes on the sample data sheets.

Reviewer's Initials/Date:

JG 5-2-88

PAGE 9 OF 13

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V  
CALIBRATION OUTLIERS  
SEMOVOLATILE HSL COMPOUNDS

(Page 1)

CASE/SAS # 9288SAS355YE

CONTRACTOR S-CUBED

Instrument # HP#	Init. Cal.	Cont. Cal.										
DATE/TIME:	4-14-88											
	RF	%RSD *	RF	%D *	RF	%D *	RF	%D *	RF	%D *	RF	%D *
Phenol												
bis(-2-Chloroethyl)Ether												
2-Chlorophenol												
1,3-Dichlorobenzene												
1,4-Dichlorobenzene												
Benzyl Alcohol												
1,2-Dichlorobenzene												
2-Methylphenol												
bis(2-chloroisopropyl)Ether												
4-Methylphenol												
N-Nitroso-Di-n-Propylamine												
Hexachloroethane												
Nitrobenzene												
Isophorone												
2-Nitrophenol												
2,4-Dimethylphenol												
Benzoic Acid												
bis(2-Chloroethoxy)Methane												
2,4-Dichlorophenol												
1,2,4-Trichlorobenzene												
Naphthalene												
4-Chloroaniline												
Hexachlorobutadiene												
4-Chloro-3-Methylphenol												
2-Methylnaphthalene												
Hexachlorocyclopentadiene												
2,4,6-Trichlorophenol												
2,4,5-Trichlorophenol												
2-Chloronaphthalene												
2-Nitroaniline												
Dimethyl Phthalate												
Acenaphthylene												
3-Nitroaniline												
Acenaphthene												
2,4-Dinitrophenol												
4-Nitrophenol												
Dibenzofuran												

LOR BLK 414  
ET831  
ET831MS  
ET831MSD  
ET832

AFFECTED  
SAMPLES:

Reviewer: AB  
Initials/Date: 5-2-88

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V  
 CALIBRATION OUTLIERS  
 SEMIVOLATILE HSL COMPOUNDS

Page 2

CASE/SAS # 9288SAS3554ECONTRACTOR S-CUBED

Instrument # HP#/	Init. Cal.	Cont. Cal.								
DATE/TIME:	9-14-88									
	RF	%RSD *	RF	%D *	RF	%D *	RF	%D *	RF	%D *
2,4-Dinitrotoluene										
2,6-Dinitrotoluene										
Diethylphthalate										
4-Chlorophenyl-phenylether										
Fluorene										
4-Nitroaniline										
4,6-Dinitro-2-Methylphenol										
N-Nitrosodiphenylamine										
4-Bromophenyl-phenylether										
Hexachlorobenzene										
Pentachlorophenol										
Phenanthrene										
Anthracene										
Di-n-Butylphthalate										
Fluoranthene										
Pyrene										
Butylbenzylphthalate										
Benzo(a)Anthracene										
bis(2-Ethylhexyl)Phthalate										
Chrysene										
Di-n-Octyl Phthalate										
Benzo(b)Fluoranthene										
Benzo(k)Fluoranthene										
Benzo(a)Pyrene										
Indeno(1,2,3-cd)Pyrene										
Dibenz(a,h)Anthracene										
Benzo(g,h,i) Perylene										

SEE PAGE 1 FOR AFFECTED SAMPLES.

\* These flags should be applied to the analytes on the sample data sheets.

Reviewer's Initials/Date:

JG 5-2-88

8/87



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION V  
 CALIBRATION OUTLIERS  
 SEMIVOLATILE HSL COMPOUNDS

CASE/SAS # 9288SAS 3554E

Page 2

CONTRACTOR

S-CUBED

Instrument #	Init. Cal.		Cont. Cal.		Cont. Cal.		Cont. Cal.		Cont. Cal.	
DATE/TIME:	RF	%RSD *	RF	%D *						
2,4-Dinitrotoluene										
2,6-Dinitrotoluene										
Diethylphthalate										
4-Chlorophenyl-phenylether										
Fluorene										
4-Nitroaniline										
4,6-Dinitro-2-Methylphenol										
N-Nitrosodiphenylamine										
4-Bromophenyl-phenylether										
Hexachlorobenzene										
Pentachlorophenol										
Phenanthrene										
Anthracene										
Di-n-Butylphthalate										
Fluoranthene										
Pyrene										
Butylbenzylphthalate										
Benzo(a)Anthracene										
bis(2-Ethylhexyl)Phthalate										
Chrysene										
Di-n-Octyl Phthalate										
Benzo(b)Fluoranthene										
Benzo(k)Fluoranthene										
Benzo(a)Pyrene										
Indeno(1,2,3-cd)Pyrene										
Dibenz(a,h)Anthracene										
Benzo(g,h,i) Perylene										

SEE PAGE 1 FOR AFFECTED SAMPLES.

\* These flags should be applied to the analytes on the sample data sheets.

Reviewer's Initials/Date: JL 5-2-88

Case: 9288-SAS 3554 E

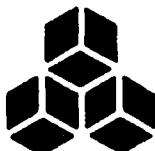
Contractor: S-CUBED

TENTATIVELY IDENTIFIED COMPOUNDS  
MATCH ASSESSMENT

NOTE: Reviewer should note directly on Organic Analysis Data Sheet (OADS) those matches that in his opinion (based on contract criteria) are unreasonable.

CRITERIA

- (1) Relative intensities of major ions (>10%) reference spectrum should be present in the sample spectrum.
- (2) Relative intensities of major ions in sample spectrum should agree to within  $\pm$  20% of reference spectrum intensities.
- (3) Molecular ions present in reference spectrum should be present in sample spectrum.
- (4) Ions present in sample spectrum, but not in reference spectrum should be reviewed for possible background contamination or presence of coeluting interferences.
- (5) Ions present in reference spectrum, but not in the sample spectrum should be reviewed for possible subtraction from the sample spectrum because of background contamination or coeluting interferences.
- (6) If, in the reviewer's opinion, no valid identification can be made the compound should be labelled as "unknown" and the initials and date of the reviewer placed on the OADS.



# S-CUBED

A Division of Maxwell Laboratories, Inc.

April 22, 1988

Volume 1 of 1

**NARRATIVE CASE 9288/SAS 3554E  
S-CUBED CONTRACT NO. 68-01-7261**

This case consists of nine waters for full analysis and one VOA only water sample. These were analyzed under the requirements of SAS 3554E which differ considerably in terms of detection limits, compound to be analyzed for as well as continuing and initial calibration requirements.

The major problem encountered was low to non-existent phenols-d<sub>5</sub> and fluorophenol surrogate recoveries in ET834, ET835, and ET836. These extracts were reinjected with similar results but reextraction could not be carried out because of the minimal sample available.

Pesticide holding times for sample extraction were missed because of a known laboratory phthalate contamination. The samples were held until laboratory corrective action showed that the extracts would not be affected by this glassware contamination.

RECORDED  
FILED  
APR 25 1988

U.S. EPA CENTRAL  
REGIONAL LAB

JoAnn Wilkinson  
JoAnn Wilkinson  
Project Manager

APR 18 1987

CENTRAL REGIONAL ENVIRONMENTAL MONITORING LABORATORY

ORGANIC INORGANIC

SUPERFUND 5011		THIS FORM IS TO BE USED FOR SAMPLES SENT TO CONTRACT ONLY	
CASE NUMBER/SAS No.		9288/3554E /37276 HADER (Ground Water) LABORATORY ANALYST DATE SHIPPED	
SUPERFUND DU NUMBER		Y905 EPA RPM or OSC (S.M.S./ICES) CONTAMINANT (STABORATOR INC.) TEST BUREAU; Ternipin (PST); SURVEYS PAGE 1 OF 1	
ACTIVITY NUMBER		C 72200	
CRL LOG NUMBER	ORGANIC TRAFFIC REPORT NUMBER	INORGANIC TRAFFIC REPORT NUMBER	
		or SAS Packing List No.	
88YLO6501	ET 831	MEW 615	X X X X X
88YLO6502	ET 832	MEW 616	X X X X X
88YLO6503	ET 833	MEW 617	X X X X X
88YLO6504	ET 834	MEW 618	X X X X X
88YLO6505	ET 835	MEW 619	X X X X X
88YLO6506	ET 836	MEW 620	X X X X X
88YLO6507	ET 837	MEW 621	X X X X X
88YLO6508	ET 838	MEW 622	X X X X X
88YLO6509	ET 839	MEW 623	X X X X X
88YLO6510	ET 840	—	X X X X X
88YLO6501	ET 841	—	X X X X X
88YLO6502	ET 842	—	X X X X X
88YLO6503	ET 843	—	X X X X X
88YLO6504	ET 844	—	X X X X X
88YLO6505	ET 845	—	X X X X X
88YLO6506	ET 846	—	X X X X X
88YLO6507	ET 847	—	X X X X X
88YLO6508	ET 848	—	X X X X X
WATER OR LIQUIDS			
ACID-BASE NEUTRAL CPDS ORGANIC SCAN UG L TOX17574			
VOLATILE ORGANIC ANALYSIS ORGANIC SCAN UG L TOX17564			
WATER POLYCHLORINATED BIPHENYLS UG L PES 17144			
WATER CHLORINATED PESTICIDES UG L PES17134			
TOTAL METALS IN WATER UG L MET111			
WATER CYANIDE UG L MIN74919			
NITRATE NITRITE MG/L MIN7284			
AMMONIA MG/L MIN7294			
RESIDUE FILTERABLE			
TDS MG/L MIN7362			
RESIDUE NON-FILT			
TSS MG/L MIN7372			
SULFATE, CHLORIDE, PH			
ALKALINITY, ACIDITY			
ACID-BASE NEUTRAL CPDS ORGANIC SCAN MG KG TOX215722			
VOLATILE ORGANIC ANALYSIS ORGANIC SCAN MG KG TOX215622			
SEDIMENTS POLYCHLORINATED BIPHENYLS MG KG PES211422			
SEDIMENT CHLORINATED PESTICIDES MG KG 211322			
TOTAL METALS MG KG MET413			
CYANIDE MG KG MIN44930			
EP TOXICITY METALS MG KG			
AMMONIA MG KG MIN42825			
SEDIMENTS OR SOILS			

# WATER SURROGATE PERCENT RECOVERY SUMMARY

Case No. 9288 (SA33554E) Contract Laboratory

S-CUBED

Contract No. 68-01-7261

SAC TRAFFIC NO.	VOLATILE						SEM-VOLATILE				PESTICIDE	
	TOLUENE-DS (100-110)	BTB (100-110)	1,2-DICHLORO- ETHANE-DS (100-110)	BENZENE-DS (100-110)	2-FLUORO- BENZENYL (100-110)	TERPENEYL- B14 (100-110)			PHENOL-DS (100-104)	2-FLUORO- PHENOL (100-100)	2,4,6-TRIBROMO- PHENOL (100-100)	CHLORFYL- CHLORDRENATE (100-100)
ET831	90	89	101	74	79	85			34	46	71	123
ET831MS	100	89	93	74	75	80			27	39	76	117
ET831MSD	99	92	88	70	75	85			38	50	83	131
ET832	89	91	101	71	78	81			30	44	71	118
ET833	107	93	99	101	99	110			36	47	102	120
ET834	102	90	104	94	93	102			0*	1*	31	120
ET835	98	89	108	68	63	77			0*	0*	15	103
ET836	96	87	99	53	52	69			1*	1*	39	91
ET837	102	94	110	70	73	81			53	57	81	109
ET838	108	97	111	75	80	93			40	48	84	126
ET839	106	94	109	51	56	66			19	23	56	126
ET840	106	93	107	NA	NA	NA			NA	NA	NA	NA
(1) LAB BLANK	94	90	105	108	107	118			66	84	94	96
(2) LAB BLANK	99	92	91	72	73	88			51	67	79	NA

\* VALUES ARE OUTSIDE OF CONTRACT REQUIRED QC LIMITS

\*\* ADVISORY LIMITS ONLY

Volatile: 0 out of 42; outside of QC limits

Semi-Volatile: 6 out of 78; outside of QC limits

Pesticides: 0 out of 12; outside of QC limits

Comments: (1) 44V2021 (VOA); LAB BLANK 4-11 (PEST); LBA07 (BNA)

(2) 4AV2021 (VOA). RRBAG (BNA)

(3) Reextraction not possible due to minimal sample volume.

# WATER MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Case No. 9288 (SAS 3554E) Contractor S-CUBED Contract No. 68-01-6868 7261

FRACTION	COMPOUND	CONC. SPIKE ADDED (µg/L)	SAMPLE RESULT	CONC. MS	% REC	CONC. MSD	% REC	RPO	QC LIMITS*	
									RPO	RECOVERY
SAMPLE NO. <u>ET 831</u>	1,1-Dichloroethene	12	0	12	100	11	92	9	14	81-145
	Trichloroethene	11	0	10	91	11	100	10	14	71-120
	Chlorobenzene	11	0	11	100	11	100	0	13	75-130
	Toluene	12	0	12	100	12	100	0	13	76-125
	Benzene	12	0	12	100	12	100	0	11	78-127
SAMPLE NO. <u>ET 831</u>	1,2,4-Trichlorobenzene	20	0	14	70	14	70	0	28	39-88
	Acenaphthene	20	0	15	75	15	75	0	31	46-118
	2,4 Dinitrotoluene	20	0	15	75	15	75	0	38	24-96
	Pyrene	20	0	16	80	17	85	6	31	26-127
	N-Nitroso-Di-n-Propylamine	20	0	15	75	14	70	7	38	41-116
	1,4-Dichlorobenzene	20	0	14	70	14	70	0	28	38-87
SAMPLE NO. <u>ET 831</u>	Pentachlorophenol	40	0	27	68	30	75	11	50	8-103
	Phenol	40	0	11	28	15	38	31	42	12-89
	2-Chlorophenol	40	0	23	58	26	65	12	40	27-123
	4-Chloro-3-Methylphenol	40	0	25	63	29	73	15	42	23-87
	4-Nitrophenol	40	0	12	30	15	38	22	50	10-80
SAMPLE NO. <u>ET 831</u>	Lindane	0.02	0	.021	105	.024	120	13	15	56-123
	Heptachlor	0.02	0	.022	110	.024	120	8.7	20	40-131
	Aldrin	0.02	0	.022	110	.023	115	4.4	22	40-120
	Dieldrin	0.05	0	.056	112	.060	120	7.0	18	52-126
	Endrin	0.05	0	.065	130*	.073	146*	12	21	56-121
	4,4'-DDT	0.05	0	.055	110	.061	122	10	27	38-127

\* ASTERISKED VALUES ARE OUTSIDE QC LIMITS.

RPO: VOA 0 out of 5: outside QC limits  
 B/N 0 out of 6: outside QC limits  
 ACID 0 out of 5: outside QC limits  
 PEST 0 out of 6: outside QC limits

RECOVERY: VOA 0 out of 10: outside QC limits  
 B/N 0 out of 12: outside QC limits  
 ACID 0 out of 10: outside QC limits  
 PEST 2 out of 12: outside QC limits

Comments: \_\_\_\_\_

\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

## **METHOD BLANK SUMMARY**

(SAS3554E)

Case No. 9288 Region 5 Contractor S-CUBED Contract No. 68-01-7267

**Contractor**

Contract No. 68-01-7261

Comments: ① ext 4/6/88

©ext 4/7/88

(A) also case holding blank

ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
: Sample Number:  
ETB31 :  
.....

Laboratory Name: S-CUBED  
Lab Sample ID No: 44V2031 (VOA)  
Sample Matrix: WATER  
Data Release Authorized By: [Signature]

Case No: 9288 (SAS3354E)  
BC Report No: N.R  
Contract No: 69-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentrations Low-Medium (Circle One)

Date Extracted/Prepared: 04-04-88

Date Analyzed: 04-04-88

Conc/Dil Factor: 1.0 pH: N.R

Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	1.3 B/U	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYL VINYLETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANDNE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYL BENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		M-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read:U=Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value:This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J) .If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/uL in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/ 3554E

Sample Number  
ET 831

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 6 Apr 88  
Date Analyzed 15 APR 88  
Conc 'Dil Factor 1.0  
Percent Moisture (Decanted) NA

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug / l or ug / Kg (Circle One)
108-95-2	Phenol	2.0 U
111-44-4	bis(2-Chloroethyl)Ether	1.5 U
95-57-8	2-Chlorophenol	2.0 U
541-73-1	1, 3-Dichlorobenzene	2.0 U
106-46-7	1, 4-Dichlorobenzene	2.0 U
100-51-6	Benzyl Alcohol	2.0 U
95-50-1	1, 2-Dichlorobenzene	2.5 U
95-48-7	2-Methylphenol	1.0 U
39638-32-8	bis(2-chloroisopropyl)Ether	2.5 U
106-44-5	4-Methylphenol	1.0 U
621-64-7	N-Nitroso-Di-n-Propylamine	1.5 U
67-72-1	Hexachloroethane	2.0 U
98-95-3	Nitrobenzene	2.5 U
78-59-1	Isophorone	2.5 U
88-75-5	2-Nitrophenol	2.0 U
105-67-9	2, 4-Dimethylphenol	2.0 U
65-85-0	Benzoic Acid	3.0 U
111-91-1	bis(2-Chloroethoxy)Methane	2.5 U
120-83-2	2, 4-Dichlorophenol	2.0 U
120-82-1	1, 2, 4-Trichlorobenzene	2.0 U
91-20-3	Naphthalene	2.0 U
106-47-8	4-Chloraniline	2.0 U
87-68-3	Hexachlorobutadiene	2.5 U
59-50-7	4-Chloro-3-Methylphenol	1.5 U
91-57-6	2-Methylnaphthalene	2.0 U
77-47-4	Hexachlorocyclopentadiene	2.0 U
88-06-2	2, 4, 6-Trichlorophenol	1.5 U
95-95-4	2, 4, 5-Trichlorophenol	1.5 U
91-58-7	2-Chloronaphthalene	1.5 U
88-74-4	2-Nitroaniline	1.0 U
131-11-3	Dimethyl Phthalate	1.5 U
208-96-8	Acenaphthylene	1.5 U
99-09-2	3-Nitroaniline	2.5 U

CAS Number		ug / l or ug / Kg (Circle One)
83-32-9	Acenaphthene	1.5 U
51-28-5	2, 4-Dinitrophenol	1.5 U
100-02-7	4-Nitrophenol	1.5 U
132-64-9	Dibenzofuran	1.0 U
121-14-2	2, 4-Dinitrotoluene	1.0 U
606-20-2	2, 6-Dinitrotoluene	1.0 U
84-66-2	Diethylphthalate	1.5
7005-72-3	4-Chlorophenyl-phenylether	1.0 U
86-73-7	Fluorene	1.0 U
100-01-6	4-Nitroaniline	3.0 U
534-52-1	4, 6-Dinitro-2-Methylphenol	1.5 U
86-30-6	N-Nitrosodiphenylamine (1)	1.5 U
101-55-3	4-Bromophenyl-phenylether	1.5 U
118-74-1	Hexachlorobenzene	1.5 U
87-86-5	Pentachlorophenol	2.0 U
85-01-8	Phenanthrene	1.0 U
120-12-7	Anthracene	2.5 U
84-74-2	Di-n-Butylphthalate	2.0 U
206-44-0	Fluoranthene	1.5 U
129-00-0	Pyrene	1.5 U
85-68-7	Butylbenzylphthalate	3.5 U
91-94-1	3, 3'-Dichlorobenzidine	NR
56-55-3	Benz(a)Anthracene *	1.5 U
117-81-7	bis(2-Ethylhexyl)Phthalate	1.9
218-01-9	Chrysene *	—
117-84-0	Di-n-Octyl Phthalate	1.5 U
205-99-2	Benz(b)Fluoranthene *	1.5 U
207-08-9	Benz(k)Fluoranthene *	—
50-32-8	Benz(a)Pyrene	2.0 U
193-39-5	Indeno(1, 2, 3-cd)Pyrene	3.5 U
53-70-3	Dibenzo(a, h)Anthracene	2.5 U
191-24-2	Benzolig. n, i)Perylene	4.0 U

(1)-Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL.

\*\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL.  
Form 1

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
ET 831

Organics Analysis Data Sheet  
(Page 3)

Pesticide/PCBs

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 4-11-88  
Date Analyzed 4-12-88  
Conc/Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-89-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxide	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
30-57-1	Dieldrin	0.010 $\mu$
72-55-9	4, 4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-85-9	Endosulfan II	0.010 $\mu$
72-54-8	4, 4'-DDD	0.020 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4, 4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53494-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53469-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
7442-93-4	DODIN ALDENEYDE	0.030 $\mu$

$V_1$  = Volume of extract injected ( $\mu$ l)

$V_2$  = Volume of water extracted (ml)

$W_3$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract ( $\mu$ l)

$v_s$  1000 ml or  $w_s$  NA  $v_1$  1.00 ml  $v_t$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Sample Number

ET831

44V2031(VOA)

ET831(BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l) (ug/kg)
1. NA	UNKNOWN	VOA	277	1.75
2.				
3. 21964498	1,13-tetradecadiene	BNA	1280	3.5
4.				
5.				
6.				
7.				
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29.				
30.				

ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
: Sample Number:  
ET832 :  
.....

Laboratory Name: S-CUBED  
Lab Sample ID No: 44V2051 (VOA)  
Sample Matrix: WATER  
Data Release Authorized By: [Signature]

Case No: 9288 (SAS3554E)  
QC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration: Low Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	0.07-0.8 U	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-66-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-13-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-9	2-CHLOROETHYLVINYLETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYLBENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		M-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnotes should read:U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J). If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/ul in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288 / 3554E

Sample Number  
ET 832

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 4-6-88  
Date Analyzed 4-15-88  
Conc.'Dil Factor 1.0  
Percent Moisture (Decanted) N/A

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	2.0 U
111-44-4	bis(2-Chloroethyl)Ether	1.5 U
95-57-8	2-Chlorophenol	2.0 U
541-73-1	1,3-Dichlorobenzene	2.0 U
106-46-7	1,4-Dichlorobenzene	2.0 U
100-51-6	Benzyl Alcohol	2.0 U
95-50-1	1,2-Dichlorobenzene	2.5 U
95-48-7	2-Methylphenol	1.0 U
39638-32-9	bis(2-chloroisopropyl)Ether	2.5 U
106-44-5	4-Methylphenol	1.0 U
621-64-7	N-Nitroso-Di-n-Propylamine	1.5 U
67-72-1	Hexachloroethane	2.0 U
98-95-3	Nitrobenzene	2.5 U
78-59-1	Isophorone	2.5 U
88-75-5	2-Nitrophenol	2.0 U
105-67-9	2,4-Dimethylphenol	2.0 U
65-85-0	Benzoic Acid	3.0 U
111-91-1	bis(2-Chloroethoxy)Methane	2.5 U
120-83-2	2,4-Dichlorophenol	2.0 U
120-82-1	1,2,4-Trichlorobenzene	2.0 U
91-20-3	Naphthalene	2.0 U
106-47-8	4-Chloroaniline	2.0 U
87-68-3	Hexachlorobutadiene	2.5 U
59-50-7	4-Chloro-3-Methylphenol	1.5 U
91-57-6	2-Methylnaphthalene	2.0 U
77-47-4	Hexachlorocyclopentadiene	2.0 U
88-06-2	2,4,6-Trichlorophenol	1.5 U
95-95-4	2,4,5-Trichlorophenol	1.5 U
91-58-7	2-Chloronaphthalene	1.5 U
88-74-4	2-Nitroaniline	1.0 U
131-11-3	Dimethyl Phthalate	1.5 U
208-96-8	Acenaphthylene	1.5 U
99-09-2	3-Nitroaniline	2.5 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	1.5 U
51-28-5	2,4-Dinitrophenol	1.5 U
100-02-7	4-Nitrophenol	1.5 U
132-84-9	Dibenzofuran	1.0 U
121-14-2	2,4-Dinitrotoluene	1.0 U
606-20-2	2,6-Dinitrotoluene	1.0 U
84-86-2	Diethylphthalate	1.0 U
7005-72-3	4-Chlorophenyl-phenylether	1.0 U
86-73-7	Fluorene	1.0 U
100-01-6	4-Nitroaniline	3.0 U
534-82-1	4,6-Dinitro-2-Methylphenol	1.5 U
86-30-6	N-Nitrosodiphenylamine (1)	1.5 U
101-55-3	4-Bromophenyl-phenylether	1.5 U
118-74-1	Hexachlorobenzene	1.5 U
87-86-5	Pentachlorophenol	2.0 U
85-01-8	Phenanthrene	1.0 U
120-12-7	Anthracene	2.5 U
84-74-2	Di-n-Butylphthalate	2.0 U
206-44-0	Fluoranthene	1.5 U
129-00-0	Pyrene	1.5 U
85-88-7	Butylbenzylphthalate	3.5 U
91-84-1	3,3'-Dichlorobenzidine	N.R.
56-55-3	Benz(a)Anthracene *	1.5 U
117-81-7	bis(2-Ethylhexyl)Phthalate	1.0 U
218-01-9	Chrysene *	—
117-84-0	Di-n-Octyl Phthalate	1.5 U
205-99-2	Benz(b)Fluoranthene * *	1.5 U
207-08-9	Benz(k)Fluoranthene * *	—
50-32-8	Benz(a)Pyrene	2.0 U
193-39-5	Indeno[1,2,3-cd]Pyrene	3.5 U
53-70-3	Dibenzo[a,h]Anthracene	2.5 U
191-24-2	Benz(g,h,i)Perylene	4.0 U

(\*)-Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL  
\*\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL  
Form 1 785

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
ET 832

Organics Analysis Data Sheet  
(Page 3)

Pesticide / PCBs

Concentration  Low      Medium (Circle One)  
Date Extracted / Prepared 04-11-88  
Date Analyzed 04-17-88  
Conc / Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug / 1g or ug / Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-89-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxyde	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
80-57-1	Dieldrin	0.010 $\mu$
72-58-9	4, 4'-DDE	0.005 $\mu$
72-20-8	Endosulfan	0.010 $\mu$
33213-65-9	Endosulfan II	0.010 $\mu$
72-54-8	4, 4'-DDD	0.020 $\mu$
1001-07-8	Endosulfan Sulfate	0.100 $\mu$
50-28-3	4, 4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53494-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Taxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53469-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
74421-93-4	ENDRIN ALDEHYDE	0.030 $\mu$

$V_1$  = Volume of extract injected ( $\mu$ l)

$V_2$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract ( $\mu$ l)

$v_s$  1000 ml or  $w_s$  NA  $v_1$  1.00 ml  $v_t$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Organics Analysis Data Sheet  
(Page 4)

Sample Number

ET832

44V2051(VOA)

ET832(BNA)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l) or ug/kg)
1.	No TIC FOUND.	VOA		
2.				
3.	No TIC Found	BNA		
4.				
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ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
Sample Number:  
ET833

Laboratory Name: S-CUBED  
Lab Sample ID No: 44V2041 (VOA)  
Sample Matrix: WATER  
Data Release Authorized By: [Signature]

Case No: 9288 (SASJ354E)  
QC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration Low Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number		<u>ug/l</u> or ug/kg (Circle One)	CAS Number		<u>ug/l</u> or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DI-BROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	1.0 0.84-J-B U	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYL VINYLETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYL BENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		M-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U flag (eg. 10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read: U=Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J). If limit of detection is 10 ug/L and a concentration of 3 ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/uL in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/3554E

Sample Number  
ET833

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted / Prepared 4-7-88  
Date Analyzed 4-10-88  
Conc 'Dil Factor 1.0  
Percent Moisture (Decanted) NA

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	2.0 U
111-44-4	bis(2-Chloroethyl)Ether	1.5 U
95-57-8	2-Chlorophenol	2.0 U
541-73-1	1, 3-Dichlorobenzene	2.0 U
106-46-7	1, 4-Dichlorobenzene	2.0 U
100-51-6	Benzyl Alcohol	2.0 U
95-50-1	1, 2-Dichlorobenzene	2.5 U
95-48-7	2-Methylphenol	1.0 U
39638-32-9	bis(2-chloroisopropyl)Ether	2.5 U
106-44-5	4-Methylphenol	1.0 U
621-64-7	N-Nitroso-Di-n-Propylamine	1.5 U
67-72-1	Hexachloroethane	2.0 U
98-95-3	Nitrobenzene	2.5 U
78-59-1	Isophorone	2.5 U
88-75-5	2-Nitrophenol	2.0 U
105-67-9	2, 4-Dimethylphenol	2.0 U
65-85-0	Benzoic Acid	3.0 U
111-91-1	bis(2-Chloroethoxy)Methane	2.5 U
120-83-2	2, 4-Dichlorophenol	2.0 U
120-82-1	1, 2, 4-Trichlorobenzene	2.0 U
91-20-3	Naphthalene	2.0 U
106-47-8	4-Chloraniline	2.0 U
87-68-3	Hexachlorobutadiene	2.5 U
59-50-7	4-Chloro-3-Methylphenol	1.5 U
91-57-6	2-Methylnaphthalene	2.0 U
77-47-4	Hexachlorocyclopentadiene	2.0 U
88-06-2	2, 4, 6-Trichlorophenol	1.5 U
95-95-4	2, 4, 5-Trichlorophenol	1.5 U
91-58-7	2-Chloronaphthalene	1.5 U
88-74-4	2-Nitroaniline	1.0 U
131-11-3	Dimethyl Phthalate	1.5 U
208-96-8	Acenaphthylene	1.5 U
99-09-2	3-Nitroaniline	2.5 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	1.5 U
51-28-5	2, 4-Dinitrophenol	1.5 U
100-02-7	4-Nitrophenol	1.5 U
132-64-9	Dibenzofuran	1.0 U
121-14-2	2, 4-Dinitrotoluene	1.0 U
606-20-2	2, 6-Dinitrotoluene	1.0 U
84-66-2	Diethylphthalate	1.0 U
7005-72-3	4-Chlorophenyl-phenylether	1.0 U
86-73-7	Fluorene	1.0 U
100-01-6	4-Nitroaniline	3.0 U
534-52-1	4, 6-Dinitro-2-Methylphenol	1.5 U
86-30-6	N-Nitrosodiphenylamine (1)	1.5 U
101-55-3	4-Bromophenyl-phenylether	1.5 U
118-74-1	Hexachlorobenzene	1.5 U
87-86-5	Pentachloropheno!	2.0 U
85-01-8	Phenanthrene	1.0 U
120-12-7	Anthracene	2.5 U
84-74-2	Di-n-Butylphthalate	2.0 U
206-44-0	Fluoranthene	1.5 U
129-00-0	Pyrene	1.5 U
35-88-7	Butylbenzylphthalate	3.5 U
91-94-1	3, 3'-Dichlorobenzidine	NR
56-55-3	Benzo(a)Anthracene *	1.5 U
117-81-7	bis(2-Ethylhexyl)Phthalate	1.0 U
218-01-9	Chrysene *	—
117-84-0	Di-n-Octyl Phthalate	1.5 U
205-99-2	Benzo(b)Fluoranthene *	1.5 U
207-08-9	Benzo(k)Fluoranthene * *	—
50-32-8	Benzo(a)Pyrene	2.0 U
193-39-5	Indeno(1, 2, 3-cd)Pyrene	3.5 U
53-70-3	Dibenzo( a, h)Anthracene	2.5 U
191-24-2	Benzo(g, h, i)Perylene	4.0 U

(1)-Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

\*\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

ET 833  
Sample Number

Organics Analysis Data Sheet  
(Page 3)

Concentration  Medium (Circle One) GPC Cleanup Dyes No  
Dose Extraced/Prepared QY-11-88 Separatory Funnel Extraction Yes  
Dose Analyzed Q4-13-88 Continuous Liquid - Liquid Extraction Yes  
Conc/Dil Factor 1.0  
Percent Moisture (decanited) NA

Number  
CAS  
Name  
Qg/g, mg/kg, %  
Available On

CAS	Name	Qg/g, mg/kg, %	Number
319-84-6	Aldrin-BHC	0.010 M	
319-85-7	Beta-BHC	0.005 M	
319-86-8	Delta-BHC	0.005 M	
319-89-9	Gamm-BHC (lindane)	0.005 M	
76-44-8	Heptachlor	0.030 M	
309-00-2	Aldrin	0.005 M	
1024-57-3	Heptachlor Epoxide	0.005 M	
359-88-8	Endosulfan I	0.010 M	
60-57-1	Dieldrin	0.010 M	
72-55-9	A, A-DDE	0.005 M	
72-30-8	Endrin	0.010 M	
33213-65-9	Endosulfan II	0.010 M	
1031-07-8	Endosulfan Sulphate	0.100 M	
50-28-3	A, A-DDT	0.020 M	
72-43-5	Methoxychlor	0.020 M	
57-74-9	Chlordane	0.020 M	
3001-35-2	Toluaphane	0.125 M	
12674-11-2	Aroclor-1016	0.10 M	
11104-28-2	Aroclor-1221	N/A	
11141-16-5	Aroclor-1232	N/A	
53469-21-9	Aroclor-1247	0.10 M	
12672-29-6	Aroclor-1248	0.10 M	
11097-69-1	Aroclor-1254	0.10 M	
11096-82-5	Aroclor-1260	0.10 M	
7621-93-4	DDT-Dieldrin	0.030 M	

Pesticide / PCBs

S-CUBED

Labatory Name 9288 (3554E) Case No.

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Sample Number

ET833

44V 2041 (VOA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (mg/l) or (ug/kg)
1.	No TIC FOUND.	VOA		
2.				
3. 21964498	1,13 tetradecadiene	BNA	1391	12 JB
4. NP	UNKNOWN		1396	6 J
5.				
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ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
: Sample Number:  
ETB34 :  
.....

Laboratory Name: S-CUBED  
Lab Sample ID No: 44V2061 (VOA)  
Sample Matrix: WATER  
Data Release Authorized By: JP

Case No: 9288 (SAS3554E)  
QC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration: Low, Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number	ug/l by ug/kg (Circle One)	CAS Number	ug/l by ug/kg (Circle One)
74-87-3 CHLOROMETHANE	10 U	78-87-5 1,2-DICHLOROPROPANE	1.5 U
74-83-9 BROMOMETHANE	10 U <u>U</u>	10061-02-6 TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4 VINYL CHLORIDE	10 U <u>U</u>	79-01-6 TRICHLOROETHENE	1.5 U
75-00-3 CHLOROETHANE	1.5 U	124-48-1 DIBROMODICHLOROMETHANE	1.5 U
75-09-2 METHYLENE CHLORIDE	1.0 -0.74 J-B U	79-00-5 1,1,2-TRICHLOROETHANE	1.5 U
67-64-1 ACETONE	75 U	71-43-2 BENZENE	1.5 U
75-15-0 CARBON DISULFIDE	3 U	10061-01-5 CIS-1,3-DICHLOROPROPENE	2 U
75-35-4 1,1-DICHLOROETHENE	1.5 U	110-75-8 2-CHLOROETHYL VINYL ETHER	1.5 U
75-34-3 1,1-DICHLOROETHANE	1.5 U	75-25-2 BROMOFORM	1.5 U
TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6 2-HEXANDNE	50 U
67-66-3 CHLOROFORM	1.5 U	108-10-1 4-METHYL-2-PENTANONE	3 U
107-06-2 1,2-DICHLOROETHANE	1.5 U	127-18-4 TETRACHLOROETHENE	1.5 U
78-93-3 2-BUTANONE	50 U	79-34-5 1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6 1,1,1-TRICHLOROETHANE	1.5 U	108-88-3 TOLUENE	1.5 U
56-23-5 CARBON TETRACHLORIDE	1.5 U	108-90-7 CHLOROBENZENE	1.5 U
108-03-4 VINYL ACETATE	15 U	100-41-4 ETHYL BENZENE	1.5 U
75-27-4 BROMODICHLOROMETHANE	1.5 U	100-42-5 STYRENE	1 U
107-02-8 ACRYLEIN	100 U	M-XYLENE	2 U
107-13-1 ACRYLONITRILE	50 U	O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Values: If the result is a value greater than or equal to the detection limit, report the value.  
 U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read:U=Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.  
 J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J) .If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

- C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/ug in the final extract should be confirmed by GC/MS.  
 B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.  
 S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/3554E

Sample Number  
ET 834

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 4-7-88  
Date Analyzed 4-10-88  
Conc 'Dil Factor 1.0  
Percent Moisture (Decanted) N/A

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	2.0 U
111-44-4	bis(2-Chloroethyl)Ether	1.5 U
95-57-8	2-Chlorophenol	2.0 U
541-73-1	1,3-Dichlorobenzene	2.0 U
106-46-7	1,4-Dichlorobenzene	2.0 U
100-51-6	Benzyl Alcohol	2.9
95-50-1	1,2-Dichlorobenzene	2.5 U
95-48-7	2-Methyphenol	1.0 U
39638-32-9	bis(2-chloroisopropyl)Ether	2.5 U
106-44-5	4-Methylphenol	1.0 U
621-64-7	N-Nitroso-Di-n-Propylamine	1.5 U
67-72-1	Hexachloroethane	2.0 U
98-95-3	Nitrobenzene	2.5 U
78-59-1	Isophorone	2.5 U
88-75-5	2-Nitrophenol	2.0 U
105-67-9	2,4-Dimethylphenol	2.0 U
65-85-0	Benzoic Acid	3.0 U
111-81-1	bis(2-Chloroethoxy)Methane	2.5 U
120-83-2	2,4-Dichlorophenol	2.0 U
120-82-1	1,2,4-Trichlorobenzene	2.0 U
91-20-3	Naphthalene	2.0 U
106-47-8	4-Chloroaniline	2.0 U
87-68-3	Hexachlorobutadiene	2.5 U
59-50-7	4-Chloro-3-Methylphenol	1.5 U
91-57-6	2-Methylnaphthalene	2.0 U
77-47-4	Hexachlorocyclopentadiene	2.0 U
88-06-2	2,4,6-Trichlorophenol	1.5 U
95-95-4	2,4,5-Trichlorophenol	1.5 U
91-58-7	2-Chloronaphthalene	1.5 U
88-74-4	2-Nitroaniline	1.0 U
131-11-3	Dimethyl Phthalate	1.5 U
208-96-8	Acenaphthylene	1.5 U
99-09-2	3-Nitroaniline	2.5 U

CAS Number	ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene 1.5 U
51-28-5	2,4-Dinitrophenol 1.5 U
100-02-7	4-Nitrophenol 1.5 U
132-64-9	Dibenzofuran 1.0 U
121-14-2	2,4-Dinitrotoluene 1.0 U
606-20-2	2,6-Dinitrotoluene 1.0 U
84-86-2	Diethylphthalate 2.9 8U
7005-72-3	4-Chlorophenyl-phenylether 1.0 U
86-73-7	Fluorene 1.0 U
100-01-8	4-Nitroaniline 3.0 U
534-52-1	4,6-Dinitro-2-Methylphenol 1.5 U
86-30-6	N-Nitrosodiphenylamine (1) 1.5 U
101-55-3	4-Bromophenyl-phenylether 1.5 U
118-74-1	Hexachlorobenzene 1.5 U
87-86-5	Pentachlorophenol 2.0 U
85-01-8	Phenanthrene 1.0 U
120-12-7	Anthracene 2.5 U
84-74-2	Di-n-Butylphthalate 2.0 U
206-44-0	Fluoranthene 1.5 U
129-00-0	Pyrene 1.5 U
85-88-7	Butylbenzylphthalate 3.5 U
91-94-1	3,3'-Dichlorobenzidine NR
56-55-3	Benz(a)Anthracene * 1.5 U
117-81-7	bis(2-Ethylhexyl)Phthalate 1.0 U
218-01-9	Chrysene *
117-84-0	Di-n-Octyl Phthalate 1.5 U
205-99-2	Benz(b)Fluoranthene ** 1.5 U
207-08-9	Benz(k)Fluoranthene **
50-32-8	Benz(a)Pyrene 2.0 U
193-39-5	Indeno(1,2,3-cd)Pyrene 3.5 U
53-70-3	Dibenz(a,h)Anthracene 2.5 U
191-24-2	Benz(o, h, i)Perylene 4.0 U

(1)-Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

\*\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

Laboratory Name S-CUBED  
Case No 9288 (3554 E)Sample Number  
ET 834Organics Analysis Data Sheet  
(Page 3)

Concentration Low Medium (Circle One)  
 Date Extracted /Prepared 04-11-88  
 Date Analyzed 04-13-88  
 Conc /Dil Factor 1.0  
 Percent Moisture (decanted) NA

GPC Cleanup  Yes  NoSeparatory Funnel Extraction  YesContinuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-89-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxide	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
60-57-1	Dieldrin	0.010 $\mu$
72-55-9	4, 4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-65-9	Endosulfan II	0.010 $\mu$
72-54-8	4, 4'-DDD	0.010 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4, 4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53494-70-5	Endrin Ketone	0.030 $\mu$
57-74-8	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53469-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
7621-93-4	ENDRIN ALDEHYDE	0.030 $\mu$

 $V_i$  = Volume of extract injected ( $\mu$ l) $V_s$  = Volume of water extracted (ml) $W_s$  = Weight of sample extracted (g) $V_t$  = Volume of total extract ( $\mu$ l) $v_s$  1000 ml or  $w_s$  NA  $v_i$  1.00 ml  $v_t$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Organics Analysis Data Sheet  
(Page 4)

Sample Number

ET834

44V2061 (VOA)

ET834 (BNA)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	No TIC FOUND.	VOA		
2.				
3. NA	UNKNOWN	BNA	243	4 J
4. NA	UNKNOWN		1226	4 J
5. 21964498	1,13 tetradecadiene		1293	11 JB
6. NA	UNKNOWN		1345	9 J
7. 1.	UNKNOWN		1358	4 J
8. NA	UNKNOWN	↓	1462	3 J
9.				
10.				
11.				
12.				
13.				
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ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
: Sample Number:  
ET835 :  
.....

Laboratory Name: S-CUBED  
Lab Sample ID No: 44V2071 (VOA), RR835(CSW)  
Sample Matrix: WATER  
Data Release Authorized By: [Signature]

Case No: 9288 (SAS3554E)  
QC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration: Low Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-3	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	/C -0.59 J-B U	79-00-3	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYL VINYLETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLORDOBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYLBENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		M-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action.(this is not necessarily the instrument detection limit.)The footnotes should read:U=Compound was analyzed for but not detected.The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value:This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J) .If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated,report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/ul in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/3554E

Organics Analysis Data Sheet  
(Page 2)

Sample Number  
E7 8.35

Concentration Low Medium (Circle One) 4 - 7 - 88  
Date Extracted / Prepared 4 - 19 - 88  
Date Analyzed 4 - 19 - 88  
Conc 'Dil Factor 1.0  
Percent Moisture (Decanted) N/A

Semivolatile Compounds

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number	ug / 1g ug / Kg (Circle One)
108-95-2 Phenol	2.0 u
111-44-4 bis(1,2-Chloroethyl)Ether	1.5 u
95-57-8 2-Chlorophenol	2.0 u
541-73-1 1,3-Dichlorobenzene	2.0 u
106-46-7 1,4-Dichlorobenzene	2.0 u
100-51-6 Benzyl Alcohol	2.0 u
95-50-1 1,2-Dichlorobenzene	2.5 u
95-48-7 2-Methylphenol	1.0 u
39638-32-8 bis(2-chloroisopropyl)Ether	2.5 u
106-44-5 4-Methylphenic	1.0 u
821-64-7 N-Nitroso-D-n-Propylamine	1.5 u
67-72-1 Hexachloroethane	3.0 u
98-95-3 Nitrobenzene	2.5 u
78-59-1 Isophorone	2.5 u
88-75-5 2-Nitrophenol	2.0 u
105-67-9 2,4-Dimethylphenol	2.0 u
65-85-0 Benzoic Acid	3.0 u
111-91-1 bis(1,2-Chloroethoxy)Methane	2.5 u
120-83-2 2,4-Dichlorophenol	2.0 u
120-82-1 1,2,4-Trichlorobutene	2.0 u
91-20-3 Nonaphthalene	2.0 u
106-47-8 4-Chloraniline	2.0 u
87-68-3 Hexachlorobutadiene	2.5 u
59-50-7 4-Chloro-3-Methylphenol	1.5 u
91-57-6 2-Methylnaphthalene	2.0 u
77-47-4 Hexachlorocyclohexadiene	2.0 u
88-06-2 2,4,6-Trichlorophenol	1.5 u
95-95-4 2,4,5-Trichlorophenol	1.5 u
91-58-7 2-Chloronaphthalene	1.5 u
88-74-4 2-Nitroaniline	1.0 u
131-11-3 Dimethyl Phthalate	1.5 u
208-96-8 Acenaphthylene	1.5 u
99-09-2 3-Nitroaniline	1.5 u

CAS Number	ug / 1g or ug / Kg (Circle One)
83-32-9 Acenaphthene	1.5 u
51-28-5 2,4-Dinitrophenol	1.5 u
100-02-7 4-Nitrophenol	1.5 u
132-84-9 Dibenzofuran	1.0 u
121-16-2 2,4-Dinitrotoluene	1.0 u
606-20-2 2,6-Dinitrotoluene	1.0 u
94-66-2 Diethylphthalate	1.0 u
7005-72-3 4-Chlorophenyl-phenylether	1.0 u
86-73-7 Fluorene	1.0 u
100-01-6 4-Nitroaniline	3.0 u
534-52-1 4,6-Dinitro-2-Methylphenol	1.5 u
96-30-6 N-Nitrosodiphenylamine (1)	1.5 u
101-55-3 4-Bromophenyl-phenylether	1.5 u
118-74-1 Hexachlorobenzene	1.5 u
87-86-5 Phenachlorophenol	2.0 u
85-01-8 Phenanthrene	1.0 u
120-12-7 Anthracene	2.5 u
84-74-2 Di-n-Butylphthalate	2.0 u
206-44-0 Fluoranthene	1.5 u
129-00-0 Pyrene	1.5 u
85-68-7 Butylbenzylphthalate	3.5 u
91-94-1 3,3'-Dichlorobenzidine	NR
56-55-3 Benz(a)Anthracene *	1.5 u
117-81-7 Bis(2-Ethylhexyl)Phthalate	1.0 u
218-01-9 Chrysene *	—
117-94-0 Di-n-Octyl Phthalate	1.5 u
205-99-2 Benzo(b)Fluoranthene **	1.5 u
207-08-9 Benzo(k)Fluoranthene **	—
50-32-8 Benzo(a)Pyrene	2.0 u
193-39-5 Indeno(1,2,3-cd)Pyrene	3.5 u
53-70-3 Dibenz(a,h)Acene	2.5 u
191-24-2 Benzo(g,h,i)Perylene	4.0 u

\*) Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL  
\*\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

Form 1

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
ET 835

Organics Analysis Data Sheet  
(Page 3)

Pesticide/PCBs

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 04-11-88  
Date Analyzed 04-17-88  
Conc 'Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/1g ug/Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-89-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxide	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
60-57-1	Dieldrin	0.010 $\mu$
72-55-9	4, 4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-65-9	Endosulfan II	0.010 $\mu$
72-54-8	4, 4'-DDD	0.020 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4, 4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53494-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53469-21-9	Aroclor-1247	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-59-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
7421-93-4	ENDRIN ALDEHYDE	0.030 $\mu$

$V_1$  = Volume of extract injected ( $\mu$ l)

$V_3$  = Volume of water extracted (ml)

$W_3$  = Weight of sample extracted (g)

$V_1$  = Volume of total extract ( $\mu$ l)

$V_3$  1000 ml or  $W_3$  NA  $V_1$  1.00 ml  $v_1$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Sample Number

ET835

44V2071(VOA)

RR835(BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l) <sup>(ug/kg)</sup> ug/kg)
1.	No TIC FOUND.	VOA		
2.				
3. NA	UNKNOWN	BNA	238	8.5
4.				
5.				
6.				
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ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
Sample Number:  
ET836 :

Laboratory Name: S-CUBED  
Lab Sample ID No: 44V2081 (VOA), RR836(ABN)  
Sample Matrix: WATER  
Data Release Authorized By: WA

Case No: 9288 (SAS3554E)  
QC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration: Low, Medium (Circle One)

Date Extracted/Prepared: 04-04-88

Date Analyzed: 04-04-88

Conc/Dil Factor: 1.0 pH: N.R

Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLORDMETHANE	10 U	78-87-3	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	1.0 0.64-0.84 U	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYLVINYLETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYLBENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		M-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Values: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg. 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnotes should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J). If limit of detection is 10 ug/L and a concentration of 3 ug/L is calculated, report as 3J.

- C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/ul in the final extract should be confirmed by GC/MS.
- B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288 / 3554E

**Organics Analysis Data Sheet**  
(Page 2)

Sample Number  
**E T 836**

Concentration Low Medium (Circle One)  GPC Cleanup  Yes  No  
Date Extracted / Prepared 4 - 7 - 88 Separatory Funnel Extraction  Yes  
Date Analyzed 4 - 19 - 88 Continuous Liquid - Liquid Extraction  Yes  
Conc 'Dil Factor 1.0  
Percent Moisture (Decanted) N/A

**Semivolatile Compounds**

CAS Number	ug / gr ug / Kg (Circle One)
108-95-2 Phenol	2.5 <input checked="" type="checkbox"/>
111-04-4 bis(2-Chloroethyl)Ether	1.5 <input checked="" type="checkbox"/>
95-57-8 2-Chlorophenol	2.0 <input checked="" type="checkbox"/>
541-73-1 1,3-Dichlorobenzene	2.0 <input checked="" type="checkbox"/>
106-46-7 1,4-Dichlorobenzene	2.0 <input checked="" type="checkbox"/>
100-51-6 Benzyl Alcohol	2.0 <input checked="" type="checkbox"/>
95-50-1 1,2-Dichlorobenzene	2.5 <input checked="" type="checkbox"/>
95-48-7 2-Methylphenol	1.0 <input checked="" type="checkbox"/>
39638-32-9 bis(2-Chloroisopropyl)Ether	2.5 <input checked="" type="checkbox"/>
106-44-5 4-Methyliphenic	1.0 <input checked="" type="checkbox"/>
621-84-7 N-Nitroso-D-n-Propylamine	1.5 <input checked="" type="checkbox"/>
67-72-1 Hexachloroethane	2.0 <input checked="" type="checkbox"/>
98-35-3 Nitrobenzene	2.5 <input checked="" type="checkbox"/>
78-59-1 Naphtalene	2.5 <input checked="" type="checkbox"/>
88-75-5 2-Nitrophenol	2.0 <input checked="" type="checkbox"/>
105-67-9 2,4-Dimethylphenol	2.0 <input checked="" type="checkbox"/>
65-85-0 Benzoic Acid	3.0 <input checked="" type="checkbox"/>
111-81-1 bis(2-Chloroethoxy)Methane	2.5 <input checked="" type="checkbox"/>
120-83-2 2,4-Dichlorophenol	2.0 <input checked="" type="checkbox"/>
120-82-1 1,2,4-Trichlorobenzene	2.0 <input checked="" type="checkbox"/>
91-20-3 Naphthalene	2.0 <input checked="" type="checkbox"/>
106-47-8 4-Chloroaniline	2.0 <input checked="" type="checkbox"/>
87-68-3 Hexachlorobutadiene	2.5 <input checked="" type="checkbox"/>
59-50-7 4-Chloro-3-Methylphenol	1.5 <input checked="" type="checkbox"/>
91-57-6 2-Methylnaphthalene	2.0 <input checked="" type="checkbox"/>
77-47-4 Hexachlorocyclohexadiene	2.0 <input checked="" type="checkbox"/>
88-06-2 2,4,6-Trichlorophenol	1.5 <input checked="" type="checkbox"/>
95-95-4 2,4,5-Trichlorophenol	1.5 <input checked="" type="checkbox"/>
91-58-7 2-Chloronaphthalene	1.5 <input checked="" type="checkbox"/>
88-74-4 2-Nitroaniline	1.0 <input checked="" type="checkbox"/>
131-11-3 Dimethyl Phthalate	1.5 <input checked="" type="checkbox"/>
208-96-8 Acenaphthylene	1.5 <input checked="" type="checkbox"/>
99-09-2 3-Nitroaniline	2.5 <input checked="" type="checkbox"/>

CAS Number	ug / gr ug / Kg (Circle One)
83-32-9 Acenaphthene	1.5 <input checked="" type="checkbox"/>
51-28-5 2,4-Dinitrophenol	1.5 <input checked="" type="checkbox"/>
100-02-7 4-Nitrophenol	1.5 <input checked="" type="checkbox"/>
132-64-9 Dibenzofuran	1.0 <input checked="" type="checkbox"/>
121-14-2 2,4-Dinitrotoluene	1.0 <input checked="" type="checkbox"/>
608-20-2 2,6-Dinitrotoluene	1.0 <input checked="" type="checkbox"/>
94-66-2 Diethylphthalate	1.0 <input checked="" type="checkbox"/>
7005-72-3 4-Chlorophenyl-phenylether	1.0 <input checked="" type="checkbox"/>
96-73-7 Fluorene	1.0 <input checked="" type="checkbox"/>
100-01-8 4-Nitroaniline	3.0 <input checked="" type="checkbox"/>
534-52-1 4,6-Dinitro-2-Methylphenol	1.5 <input checked="" type="checkbox"/>
96-30-6 N-Nitrosodiphenylamine (11)	1.5 <input checked="" type="checkbox"/>
101-55-3 4-Bromophenyl-phenylether	1.5 <input checked="" type="checkbox"/>
118-74-1 Hexachlorobenzene	1.5 <input checked="" type="checkbox"/>
87-88-5 Pentachlorophenol	2.0 <input checked="" type="checkbox"/>
95-01-8 Phenanthrene	1.0 <input checked="" type="checkbox"/>
120-12-7 Anthracene	2.5 <input checked="" type="checkbox"/>
94-74-2 Di-n-Butylphthalate	2.0 <input checked="" type="checkbox"/>
206-44-0 Fluoranthene	1.5 <input checked="" type="checkbox"/>
129-00-0 Pyrene	1.5 <input checked="" type="checkbox"/>
95-68-7 Butylbenzylphthalate	3.5 <input checked="" type="checkbox"/>
91-94-1 3,3'-Dichlorobenzidine	NR
56-55-3 BenzeneAnthracene *	1.5 <input checked="" type="checkbox"/>
117-81-7 bis(2-Ethylhexyl)Phthalate	1.0 <input checked="" type="checkbox"/>
218-01-9 Chrysene *	—
117-84-0 D-n-Octyl Phthalate	1.5 <input checked="" type="checkbox"/>
205-98-2 Benzofluoranthene *	1.5 <input checked="" type="checkbox"/>
207-08-9 Benzotifluoranthene *	—
50-32-8 Benzofluorene	2.0 <input checked="" type="checkbox"/>
193-39-5 Indenol, 2,3-eciprene	2.5 <input checked="" type="checkbox"/>
53-70-3 Dibenz[e,h]Anthracene	2.5 <input checked="" type="checkbox"/>
191-24-2 Benzog. h. iperylene	4.0 <input checked="" type="checkbox"/>

(11) Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

# THESE TWO PARAMETERS ARE REPORTED AS A TOTAL FORM I

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
ET 836

Organics Analysis Data Sheet  
(Page 3)

Pesticide / PCBs

Concentration Low Medium (Circle One)  
Date Extracted / Prepared 04-11-88  
Date Analyzed 04-13-88  
Conc / Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug / 1 g ug / Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-89-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxide	0.005 $\mu$
959-38-8	Endosulfan I	0.010 $\mu$
60-57-1	Dieldrin	0.010 $\mu$
72-55-9	4,4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-85-9	Endosulfan II	0.010 $\mu$
72-54-8	4,4'-DDD	0.020 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4,4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53494-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53489-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
7421-93-4	ENDRIN ALDEHYDE	0.030 $\mu$

$V_1$  = Volume of extract injected ( $\mu$ l)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract ( $\mu$ l)

$v_s$  1000 ml or  $w_s$  NA  $v_1$  1.00 ml  $v_t$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Sample Number

ET 836

44V2081 (VOA)  
RR 836 (BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l) or ug/kg)
1.	No TIC FOUND.	VOA		
2.				
3. <u>NR</u>	UNKNOWN	BNA	227	4J
4. <u>21964498</u>	1,13-tetrademadiene.	BNA	1274	20 JB
5.				
6.				
7.				
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28.				
29.				
30.				

ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
Sample Number:  
ET837 :  
.....

Laboratory Name: S-CUBED  
Lab Sample ID No: 4AV2031 (VOA)  
Sample Matrix: WATER  
Data Release Authorized By: [Signature]

Case No: 9288 (SAS3554E)  
GC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentrations: Low-Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number	ug/l or ug/kg (Circle One)	CAS Number	ug/l or ug/kg (Circle One)
74-87-3 CHLOROMETHANE	10 U	78-87-5 1,2-DICHLOROPROPANE	1.5 U
74-83-9 BROMOMETHANE	10 U	10061-02-6 TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4 VINYL CHLORIDE	10 U	79-01-6 TRICHLOROETHENE	1.5 U
75-00-3 CHLOROETHANE	1.5 U	124-48-1 DIBROMOCHLOROMETHANE	1.5 U
75-09-2 METHYLENE CHLORIDE	1 8U	79-00-5 1,1,2-TRICHLOROETHANE	1.5 U
67-64-1 ACETONE	75 U	71-43-2 BENZENE	1.5 U
75-15-0 CARBON DISULFIDE	3 U	10061-01-5 CIS-1,3-DICHLOROPROPENE	2 U
75-35-4 1,1-DICHLOROETHENE	1.5 U	110-75-8 2-CHLOROETHYL VINYL ETHER	1.5 U
75-34-3 1,1-DICHLOROETHANE	1.5 U	75-29-2 BROMOFORM	1.5 U
TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6 2-HEXANONE	50 U
67-66-3 CHLOROFORM	1.5 U	108-10-1 4-METHYL-2-PENTANONE	3 U
107-06-2 1,2-DICHLOROETHANE	1.5 U	127-18-4 TETRACHLOROETHENE	1.5 U
78-93-3 2-BUTANONE	50 U	79-34-5 1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6 1,1,1-TRICHLOROETHANE	1.5 U	108-88-3 TOLUENE	1.5 U
56-23-5 CARBON TETRACHLORIDE	1.5 U	108-90-7 CHLOROBENZENE	1.5 U
108-05-4 VINYL ACETATE	15 U	100-41-4 ETHYLBENZENE	1.5 U
75-27-4 BROMODICHLOROMETHANE	1.5 U	100-42-5 STYRENE	1 U
107-02-8 ACROLEIN	100 U	N-XYLENE	2 U
107-13-1 ACRYLONITRILE	50 U	O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read:U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J) . If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/ul in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/3554E

Organics Analysis Data Sheet  
(Page 2)

Sample Number  
ET 837

Semivolatile Compounds

Concentration Low Medium (Circle One) GPC Cleanup  Yes  No  
Date Extracted / Prepared 4-7-88 Separatory Funnel Extraction  Yes  
Date Analyzed 4-14-88 Continuous Liquid - Liquid Extraction  Yes  
Cont 'Dil Factor' 1.0.  
Percent Moisture (Decanted) N/A

CAS Number	ug / ml or ug / kg (Circle One)	CAS Number	ug / gr ug / kg (Circle One)
108-95-2 Phenol	2.6	83-32-9 Acenaphthene	1.5 u
111-44-4 bis(2-Chloroethyl)Ether	1.5 u	51-28-5 2,4-Dinitrophenol	1.5 u
95-57-8 2-Chlorophenol	3.0 u	100-02-7 4-Nitrophenol	1.5 u
541-73-1 1,3-Dichlorobenzene	2.0 u	132-64-9 Dibenzofuran	1.0 u
106-46-7 1,4-Dichlorobenzene	2.0 u	121-14-2 2,4-Dinitrotoluene	1.0 u
100-51-6 Benzyl Alcohol	2.0 u	606-20-2 2,6-Dinitrotoluene	1.0 u
95-50-1 1,2-Dichlorobenzene	2.5 u	94-86-2 Diethylphthalate	1.9 <sup>11</sup> u
95-48-7 2-Methylphenol	1.0 u	7005-72-3 4-Chlorophenyl-phenylether	1.0 u
39638-32-9 bis(2-chloroethyl)ether	2.5 u	96-73-7 Fluorene	1.0 u
106-44-5 4-Methylphenol	1.0 u	100-01-6 4-Nitroaniline	3.0 u
621-64-7 N-Nitroso-D-n-Propylamine	1.5 u	534-62-1 4,6-Dinitro-2-Methylphenol	1.5 u
67-72-1 Hexachlorobutane	2.0 u	86-30-6 N-Nitrosodiphenylamine (1)	1.5 u
98-95-3 Nitrobenzene	2.5 u	101-55-3 4-Bromophenyl-phenylether	1.5 u
78-59-1 Isophorone	2.5 u	118-74-1 Hexachlorobenzene	1.5 u
88-75-5 2-Nitrophenol	2.0 u	87-86-5 Pentachlorophenol	2.0 u
105-67-9 2,4-Dimethylphenol	2.0 u	95-01-8 Phenanthrene	1.0 u
65-85-0 Benzoic Acid	3.0 u	120-12-7 Anthracene	2.5 u
111-91-1 bis(2-Chloroethyl)methane	2.5 u	94-74-2 Di-n-Butylphthalate	3.0 u
120-83-2 2,4-Dichlorophenol	2.0 u	206-44-0 Fluoranthene	1.5 u
120-82-1 1,2,4-Trichlorobenzene	2.0 u	129-00-0 Pyrene	1.5 u
91-20-3 Naphthalene	2.0 u	95-68-7 Butylbenzylphthalate	6.6 u
106-47-8 4-Chloroaniline	2.0 u	91-94-1 3,3-Dichlorobenzidine	N.R.
87-68-3 Hexachlorobutadiene	2.5 u	56-55-3 Benzene/Anthracene *	1.5 u
59-50-7 4-Chloro-3-Methylphenol	1.5 u	117-81-7 bis(2-Ethylvinyl)phthalate	1.0 u
91-57-6 2-Methylnaphthalene	2.0 u	218-01-9 Chrysene *	—
77-47-4 Hexachlorocyclohexadiene	2.0 u	117-84-0 D-n-Octyl Phthalate	1.5 u
88-06-2 2,4,6-Trichlorophenol	1.5 u	205-99-2 Benzobifluoranthene *	1.5 u
95-95-4 2,4,5-Trichlorophenol	1.5 u	207-08-9 Benzalklyfluoranthene *	—
91-58-7 2-Chloronaphthalene	1.5 u	50-32-8 Benzaldehyde	2.0 u
88-74-4 2-Nitroaniline	1.0 u	183-39-5 Indanol, 2,3-diPvrene	3.5 u
131-11-3 Dimethyl Phthalate	1.5 u	53-70-3 Dibenzyl, n-Phthalicene	2.5 u
208-98-8 Acenaphthylene	1.5 u	191-24-2 Benzalq. n. Perylene	4.0 u
99-09-2 3-Nitroaniline	2.5 u		

CAS Number	ug / ml or ug / kg (Circle One)	CAS Number	ug / gr ug / kg (Circle One)
83-32-9 Acenaphthene	1.5 u	51-28-5 2,4-Dinitrophenol	1.5 u
51-28-5 2,4-Dinitrophenol	1.5 u	100-02-7 4-Nitrophenol	1.5 u
100-02-7 4-Nitrophenol	1.5 u	132-64-9 Dibenzofuran	1.0 u
132-64-9 Dibenzofuran	1.0 u	121-14-2 2,4-Dinitrotoluene	1.0 u
121-14-2 2,4-Dinitrotoluene	1.0 u	606-20-2 2,6-Dinitrotoluene	1.0 u
606-20-2 2,6-Dinitrotoluene	1.0 u	94-86-2 Diethylphthalate	1.9 <sup>11</sup> u
94-86-2 Diethylphthalate	1.9 <sup>11</sup> u	7005-72-3 4-Chlorophenyl-phenylether	1.0 u
7005-72-3 4-Chlorophenyl-phenylether	1.0 u	96-73-7 Fluorene	1.0 u
96-73-7 Fluorene	1.0 u	100-01-6 4-Nitroaniline	3.0 u
100-01-6 4-Nitroaniline	3.0 u	534-62-1 4,6-Dinitro-2-Methylphenol	1.5 u
534-62-1 4,6-Dinitro-2-Methylphenol	1.5 u	86-30-6 N-Nitrosodiphenylamine (1)	1.5 u
86-30-6 N-Nitrosodiphenylamine (1)	1.5 u	101-55-3 4-Bromophenyl-phenylether	1.5 u
101-55-3 4-Bromophenyl-phenylether	1.5 u	118-74-1 Hexachlorobenzene	1.5 u
118-74-1 Hexachlorobenzene	1.5 u	87-86-5 Pentachlorophenol	2.0 u
87-86-5 Pentachlorophenol	2.0 u	95-01-8 Phenanthrene	1.0 u
95-01-8 Phenanthrene	1.0 u	120-12-7 Anthracene	2.5 u
120-12-7 Anthracene	2.5 u	94-74-2 Di-n-Butylphthalate	3.0 u
94-74-2 Di-n-Butylphthalate	3.0 u	206-44-0 Fluoranthene	1.5 u
206-44-0 Fluoranthene	1.5 u	129-00-0 Pyrene	1.5 u
129-00-0 Pyrene	1.5 u	95-68-7 Butylbenzylphthalate	6.6 u
95-68-7 Butylbenzylphthalate	6.6 u	91-94-1 3,3-Dichlorobenzidine	N.R.
91-94-1 3,3-Dichlorobenzidine	N.R.	56-55-3 Benzene/Anthracene *	1.5 u
56-55-3 Benzene/Anthracene *	1.5 u	117-81-7 bis(2-Ethylvinyl)phthalate	1.0 u
117-81-7 bis(2-Ethylvinyl)phthalate	1.0 u	218-01-9 Chrysene *	—
218-01-9 Chrysene *	—	117-84-0 D-n-Octyl Phthalate	1.5 u
117-84-0 D-n-Octyl Phthalate	1.5 u	205-99-2 Benzobifluoranthene *	1.5 u
205-99-2 Benzobifluoranthene *	1.5 u	207-08-9 Benzalklyfluoranthene *	—
207-08-9 Benzalklyfluoranthene *	—	50-32-8 Benzaldehyde	2.0 u
50-32-8 Benzaldehyde	2.0 u	183-39-5 Indanol, 2,3-diPvrene	3.5 u
183-39-5 Indanol, 2,3-diPvrene	3.5 u	53-70-3 Dibenzyl, n-Phthalicene	2.5 u
53-70-3 Dibenzyl, n-Phthalicene	2.5 u	191-24-2 Benzalq. n. Perylene	4.0 u
191-24-2 Benzalq. n. Perylene	4.0 u		

(1)-Cannot be separated from diphenylamine

# THREE TWO PARAMETERS ARE REPORTED AS A TOTAL

\* \* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL  
form,

7 85

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
ET 837

Organics Analysis Data Sheet  
(Page 3)

Pesticide/PCBs

Concentration Low Medium (Circle One)  
Date Extracted /Prepared 04-11-88  
Date Analyzed 04-13-88  
Conc /Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/1g or ug/Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-89-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxyde	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
50-57-1	Dieldrin	0.010 $\mu$
72-55-9	4,4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-85-9	Endosulfan II	0.010 $\mu$
72-54-8	4,4'-DDD	0.020 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4,4'-DDT	0.020 $\mu$
72-43-8	Methoxychlor	0.020 $\mu$
53494-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53469-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
7421-93-4	ENDRIN ALDEHYDE	0.030 $\mu$

$V_1$  = Volume of extract injected ( $\mu$ l)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract ( $\mu$ l)

$v_s$  1000 ml or  $w_s$  NA  $v_1$  1.00 ml  $v_t$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No. 9288 (SAS5554E)

Sample Number

ET 837

4AV2031(VOA)

ET 837(BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l) or ug/kg/l
1.	NO TIC FOUND.	VOA		
2.				
3. 31964498	1,13 tetradecadiene	BNA	1284	7 JB
4.				
5.				
6.				
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28.				
29.				
30.				

ORGANICS ANALYSIS DATA SHEET  
(Page 1)

Laboratory Name: S-CUBED  
Lab Sample ID No: V244011 (VOA)  
Sample Matrix: WATER  
Data Release Authorized By: [initials]

Case No: 9288 (SAS3554E)  
GC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration: Low Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	1 8 U	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYL VINYL ETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 U 0.30-3 μu	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYL BENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		N-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read: U=Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J). If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/uL in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/3554E

### Organics Analysis Data Sheet (Page 2)

#### Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted / Prepared 4-7-88  
Date Analyzed 4-14-88  
Conc 'Dil Factor 1.0  
Percent Moisture (Decanted) N/A

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number	ug / 1g (Circle One)	ug / 1g ug / Kg (Circle One)
108-95-2 Phenol	2.0 u	63.32.9 Acenaphthene
111-44-4 bis(2-Chloroethyl)Ether	1.5 u	61.28.5 2, 4-Dinitrophenol
95-57-8 2-Chlorophenol	2.0 u	100-02.7 4-Nitrophenol
541-73-1 1, 3-Dichlorobenzene	2.0 u	132-84.9 Dibenzofuran
106-46-7 1, 4-Dichlorobenzene	2.0 u	121-14.2 2, 4-Dinitrotoluene
100-51-6 Benzyl Alcohol	2.0 u	606-20.2 2, 6-Dinitrotoluene
95-50-1 1, 2-Dichlorobenzene	2.5 u	84-66-2 Diethylphthalate
95-48-7 2-Methylphenol	1.0 u	7005-72-3 4-Chlorophenyl-phenylether
39638-32-9 bis(2-chloroisopropyl)Ether	2.5 u	86-73-7 Fluorene
106-44-5 4-Methylphenol	1.0 u	100-01-8 4-Nitroaniline
621-84-7 N-Nitroso-D-n-Propylamine	1.5 u	534-62-1 4, 6-Dinitro-2-Methylphenol
67-72-1 Hexachloroethane	2.0 u	96-30-6 N-Nitroso-diphenylamine (11)
98-95-3 Nitrobenzene	2.5 u	101-65-3 4-Bromophenyl-phenylether
78-59-1 Isophorone	2.5 u	118-74-1 Hexachlorobenzene
88-75-5 2-Nitrophenol	2.0 u	97-96-5 Pentachlorophenol
105-67-9 2, 4-Dimethylphenol	2.0 u	85-01-8 Phenanthrene
65-85-0 Benzoic Acid	3.0 u	120-12-7 Anthracene
111-91-1 bis(2-Chloroethyl)Methane	2.5 u	84-74-2 Di-n-Butylphthalate
120-83-2 2, 4-Dichlorophenol	2.0 u	206-44-0 Fluoranthene
120-82-1 1, 2, 4-Trichlorobenzene	2.0 u	129-00-0 Pyrene
91-20-3 Naphthalene	2.0 u	85-68-7 Butylbenzylphthalate
106-47-8 4-Chloroaniline	2.0 u	91-94-1 3, 3-Dichlorobenzidine
87-68-3 Methylchlorobutadiene	2.5 u	56-55-3 Benzodihydrofuran
59-50-7 4-Chloro-3-Methylphenol	1.5 u	117-91-7 bis(2-Ethylhexyl)Phthalate
91-57-6 2-Methylnaphthalene	2.0 u	218-01-9 Chrysene *
77-47-4 Methylchlorocyclohexadiene	2.0 u	117-84-0 Di-n-Octyl Phthalate
88-06-2 2, 4, 6-Trichlorophenol	1.5 u	205-99-2 Benzodifluoranthene **
85-95-4 2, 4, 5-Trichlorophenol	1.5 u	207-08-9 Benzofluoranthene ***
91-58-7 2-Chloronaphthalene	1.5 u	50-32-8 BenzoflPyrene
88-74-4 2-Nitroaniline	1.0 u	193-39-5 Indeno[1, 2, 3-cd]Pyrene
131-11-3 Dimethyl Phthalate	1.5 u	53-70-3 Dibenz[a,h]Anthracene
208-96-8 Acenaphthylene	1.5 u	191-24-2 Benzog. h. Perylene
99-09-2 3-Nitroaniline	2.5 u	

CAS Number	ug / 1g (Circle One)	ug / 1g ug / Kg (Circle One)
63-32-9 Acenaphthene	1.5 u	
61-28-5 2, 4-Dinitrophenol	1.5 u	
100-02-7 4-Nitrophenol	1.5 u	
132-84-9 Dibenzofuran	1.0 u	
121-14-2 2, 4-Dinitrotoluene	1.0 u	
606-20-2 2, 6-Dinitrotoluene	1.0 u	
84-66-2 Diethylphthalate	1.0 u	
7005-72-3 4-Chlorophenyl-phenylether	1.0 u	
86-73-7 Fluorene	1.0 u	
100-01-8 4-Nitroaniline	2.0 u	
534-62-1 4, 6-Dinitro-2-Methylphenol	1.5 u	
96-30-6 N-Nitroso-diphenylamine (11)	1.5 u	
101-65-3 4-Bromophenyl-phenylether	1.5 u	
118-74-1 Hexachlorobenzene	1.5 u	
97-96-5 Pentachlorophenol	2.0 u	
85-01-8 Phenanthrene	1.0 u	
120-12-7 Anthracene	2.5 u	
84-74-2 Di-n-Butylphthalate	2.0 u	
206-44-0 Fluoranthene	1.5 u	
129-00-0 Pyrene	1.5 u	
85-68-7 Butylbenzylphthalate	3.5 u	
91-94-1 3, 3-Dichlorobenzidine	N.R.	
56-55-3 Benzodihydrofuran	1.5 u	
117-91-7 bis(2-Ethylhexyl)Phthalate	1.0 u	
218-01-9 Chrysene *	—	
117-84-0 Di-n-Octyl Phthalate	1.5 u	
205-99-2 Benzodifluoranthene **	1.5 u	
207-08-9 Benzofluoranthene ***	—	
50-32-8 BenzoflPyrene	2.0 u	
193-39-5 Indeno[1, 2, 3-cd]Pyrene	2.5 u	
53-70-3 Dibenz[a,h]Anthracene	2.5 u	
191-24-2 Benzog. h. Perylene	4.0 u	

(1)-Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL  
\*\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL  
form!

7 85

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
ET 838

Organics Analysis Data Sheet  
(Page 3)

Pesticide / PCBs

Concentration  Low      Medium      (Circle One)  
Date Extracted /Prepared 04-11-88  
Date Analyzed 04-13-88  
Conc / Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-89-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxide	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
80-57-1	Dieldrin	0.010 $\mu$
72-55-9	4, 4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-65-9	Endosulfan II	0.010 $\mu$
72-54-8	4, 4'-DDD	0.020 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4, 4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53484-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53489-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
7621-93-4	ENDRIN ALDENEYES	0.030 $\mu$

$V_1$  = Volume of extract injected ( $\mu$ l)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract ( $\mu$ l)

$v_s$  1000 ml or  $w_s$  NA  $v_1$  1.00 ml  $v_t$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No 9288 (SAS 3554 E)

Sample Number

ET 838

V244011 (VOA)  
ET 838 (BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration ng/Pg ug/kg)
1.	NO TIC FOUND.	VOA		
2.				
3. 21964498	1,13 tetradecadiene	BNA	1287	12 JB
4. 19812647	1,14 tetra decane diol		1290	3 JB
5.				
6.				
7.				
8.				
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ORGANICS ANALYSIS DATA SHEET  
(Page 1)

Laboratory Name: S-CUBED  
 Lab Sample ID No: V244021 (VON), R2839 (ABN)  
 Sample Matrix: WATER  
 Data Release Authorized By: [Signature]

Case No: 9288 (SAS3554E)  
 QC Report No: N.R  
 Contract No: 68-01-7261  
 Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration/Low Medium (Circle One)

Date Extracted/Prepared: 04-04-88  
 Date Analyzed: 04-04-88  
 Conc/Dil Factor: 1.0 pH: N.R  
 Percent Moisture (Not Decanted): N.R

CAS Number		<u>ug/l or ug/kg</u> (Circle One)	CAS Number		<u>ug/l or ug/kg</u> (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DI(BROMOCHLOROMETHANE)	1.5 U
75-09-2	METHYLENE CHLORIDE	1.4 <del>8U</del>	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-73-8	2-CHLOROETHYL VINYL ETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 <del>0.23-0.5U</del>	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYL BENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		N-KYLINE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg. 10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read: U=Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J). If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/µl in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/3554E

Sample Number  
ET 839

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 4-7-88  
Date Analyzed 4-19-88  
Conc.'Dil Factor 1.0  
Percent Moisture (Decanted) N/A

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug /Kg (Circle One)
108-95-2	Phenol	2.0 U
111-44-4	bis(2-Chloroethyl)Ether	1.5 U
95-57-8	2-Chlorophenol	2.0 U
541-73-1	1, 3-Dichlorobenzene	2.0 U
106-46-7	1, 4-Dichlorobenzene	2.0 U
100-51-6	Benzyl Alcohol	2.0 U
95-50-1	1, 2-Dichlorobenzene	2.5 U
95-48-7	2-Methylphenol	1.0 U
3963B-32-9	bis(2-chloroisooxypropyl)Ether	2.5 U
106-44-5	4-Methylphenol	1.0 U
621-64-7	N-Nitroso-Di-n-Propylamine	1.5 U
67-72-1	Hexachloroethane	2.0 U
98-95-3	Nitrobenzene	2.5 U
78-59-1	Isophorone	2.5 U
88-75-5	2-Nitrophenol	2.0 U
105-67-9	2, 4-Dimethylphenol	2.0 U
65-85-0	Benzoic Acid	3.0 U
111-91-1	bis(2-Chloroethoxy)Methane	2.5 U
120-83-2	2, 4-Dichlorophenol	2.0 U
120-82-1	1, 2, 4-Trichlorobenzene	2.0 U
91-20-3	Naphthalene	2.0 U
106-47-8	4-Chloroaniline	2.0 U
87-68-3	Hexachlorobutadiene	2.5 U
59-50-7	4-Chloro-3-Methylphenol	1.5 U
91-57-6	2-Methylnaphthalene	2.0 U
77-47-4	Hexachlorocyclopentadiene	2.0 U
88-06-2	2, 4, 6-Trichlorophenol	1.5 U
95-95-4	2, 4, 5-Trichlorophenol	1.5 U
91-58-7	2-Chloronaphthalene	1.5 U
88-74-4	2-Nitroaniline	1.0 U
131-11-3	Dimethyl Phthalate	1.5 U
208-96-8	Acenaphthylene	1.5 U
89-09-2	3-Nitroaniline	2.5 U

CAS Number		ug/l or ug /Kg (Circle One)
83-32-9	Acenaphthene	1.5 U
51-28-5	2, 4-Dinitrophenol	1.5 U
100-02-7	4-Nitrophenol	1.5 U
132-64-9	Dibenzofuran	1.0 U
121-14-2	2, 4-Dinitrotoluene	1.0 U
606-20-2	2, 6-Dinitrotoluene	1.0 U
84-66-2	Diethylphthalate	1.2 V B
7005-72-3	4-Chlorophenyl-phenylether	1.0 U
86-73-7	Fluorene	1.0 U
100-01-8	4-Nitroaniline	3.0 U
534-52-1	4, 6-Dinitro-2-Methylphenol	1.5 U
86-30-6	N-Nitrosodiphenylamine (1)	1.5 U
101-55-3	4-Bromophenyl-phenylether	1.5 U
118-74-1	Hexachlorobenzene	1.5 U
87-86-5	Pentachloropheno!	2.0 U
85-01-8	Phenanthrene	1.0 U
120-12-7	Anthracene	2.5 U
84-74-2	Di-n-Butylphthalate	2.0 U
206-44-0	Fluoranthene	1.5 U
129-00-0	Pyrene	1.5 U
85-68-7	Butylbenzylphthalate	7.1
91-94-1	3, 3'-Dichlorobenzidine	NR
56-55-3	Benz(a)Anthracene *	1.5 U
117-81-7	bis(2-Ethylhexyl)Phthalate	1.0 U
218-01-9	Chrysene *	—
117-84-0	Di-n-Octyl Phthalate	1.5 U
205-99-2	Benz(b)Fluoranthene * *	1.5 U
207-08-9	Benz(k)Fluoranthene * *	—
50-32-8	Benz(a)Pyrene	2.0 U
193-39-5	Indeno[1, 2, 3-cd]Pyrene	3.5 U
63-70-3	Dibenzo[a, h]Anthracene	2.5 U
191-24-2	Benz(g, h, i)Perylene	4.0 U

(1)-Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

\*\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
ET 839

Organics Analysis Data Sheet  
(Page 3)

Pesticide/PCBs

Concentration Low Medium (Circle One)  
Date Extracted /Prepared 04-11-88  
Date Analyzed 04-13-88  
Conc /Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-99-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxide	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
60-57-1	Dieldrin	0.010 $\mu$
72-55-9	4,4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-65-9	Endosulfan II	0.010 $\mu$
72-84-8	4,4'-DDD	0.020 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4,4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53484-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53489-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
7421-93-4	ENDRIN ALDEHYDE	0.030 $\mu$

$V_1$  = Volume of extract injected (ml)

$V_2$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ml)

$v_s$  1000 ml or  $w_s$  NA  $v_t$  1.00 ml  $v_i$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No. 9288 (SAS3554E)

Sample Number

ET839

V244021 (VOA)  
RR 839 (BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l) (ug/kg)
1.	NO TIC FOUND.	VOA		
2.				
3. <u>81964498</u>	1,13 Tetradecadiene	BNA	1274	4 JB
4. <u>NA</u>	UNKNOWN	↓	1379	8 J
5.				
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ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
: Sample Number:  
ET840 :  
.....

Laboratory Name: S-CUBED  
Lab Sample ID No: V244031 (VOA)  
Sample Matrix: WATER  
Data Release Authorized By: [initials]

Case No: 9288 (SAS3554E)  
QC Report No: N.R  
Contract No: 68-01-7261  
Date Sample Received: 04-02-88

VOLATILE COMPOUNDS

Concentration (Low) Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	1.3 U	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYL VINYL ETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANDNE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U.
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYL BENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		N-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read:U=Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J) . If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/uL in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Instrument Name 3-CURD  
Case No 9288 (SASSSE)

Organics Analysis Data Sheet  
(Page 4)

Sample Number  
ET 840

V244031 (VOA)

Tentatively Identified Compounds

CAS Number	Compound Name	Precision	RT or Scan Number	Estimated Concentration (ppm up to 100)
1	No Tic FOUND.	VOA		
2	NO ABN ANALYSIS			
3				
4				
5				
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UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

ESJ/Central Regional Laboratory  
DATA TRACKING FORM FOR CONTRACT SAMPLES

CRL Data Set No. SF 5011 CERCLIS No.   

SMO Case No. 9288 SAS 3554E Site Name and Location: HADER GROUND WATER

Name of Contractor or EPA Laboratory: S-CUBED Data User: MPCA

No. of Samples: 10 Date Samples or Data Received: 4-25-88

1. Have chain-of-custody records been received? YES  NO
  2. Have Traffic Reports or packing lists been received? YES  NO
  3. If no, are Traffic Report or packing list numbers written on the chain-of-custody record? YES  NO
  4. If no, which Traffic report or packing list numbers are missing?
- \_\_\_\_\_
- \_\_\_\_\_
- \_\_\_\_\_

Are basic data forms in? YES  NO

Number of samples claimed: 10 Number of samples received: 10

Checked by: Erlinda Luz M. Arresta Date: 4-25-88

Received by Contract Project Management Section: P.Chanilla Date: 4-25-88

Review Started: 5-2-88 Reviewer Signature: Jeanne R. Maynard

Total time spent on review: 7hr Date review completed: 5-2-88

Copied (xeroxed) by: \_\_\_\_\_ Date: \_\_\_\_\_

Mailed to Data User by: Wanda Leeman Date: 56-88

DATA USERS:

Please fill in the blanks below and return this form to: Sylvia Griffin, Data Management Coordinator, Region V, SSCRL

Data received by: \_\_\_\_\_ Date: \_\_\_\_\_

Q.A. review received by: \_\_\_\_\_ Date: \_\_\_\_\_

Inorganic Data Complete [ ]. Suitable for Intended Purposes [ ]  [ ] if acceptable.  
Organic Data Complete [ ]. Suitable for Intended Purposes [ ] List problems below.  
Dioxin Data Complete [ ]. Suitable for Intended Purposes [ ]  
SAS Data Complete [ ]. Suitable for Intended Purposes [ ]

See Attached "Missing Data Request Form" [ ]  
PROBLEMS: Please indicate reasons (if any) why data are not suitable for your uses.  
Other problems.

\_\_\_\_\_

\_\_\_\_\_

Received by Data Management Coordinator, CRL for File: Date: \_\_\_\_\_

Signature: \_\_\_\_\_

ORGANICS ANALYSIS DATA SHEET  
(Page 1)

.....  
Sample Number:  
LAB BLANK:

Laboratory Name: S-CUBED  
 Lab Sample ID No: 44Y2021 (VOA), RR8AG (ABW)  
 Sample Matrix: WATER Lab blank 9-11 (pre)  
 Data Release Authorized By: [initials]

Case No: 9288 (SAS3554E)  
 QC Report No: N.R  
 Contract No: 68-01-7261  
 Date Sample Received: N.R

VOLATILE COMPOUNDS

Concentration  Medium (Circle One)

Date Extracted/Prepared: 04-04-88

Date Analyzed: 04-04-88

Conc/Dil Factor: 1.0 pH: N.R

Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	1.3	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYL VINYLETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-55-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYL BENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		M-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

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 J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J) . If limit of detection is 10 ug/L and a concentration of 3 ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/ul in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 92 89/3554E

Sample Number  
Lab Blank 4-6

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 4-6-88  
Date Analyzed 4-13-88  
Conc.'Dil Factor 1.0  
Percent Moisture (Decanted) N/A

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	2.0 U
111-44-4	bis(2-Chloroethyl)Ether	1.5 U
95-57-8	2-Chlorophenol	2.0 U
541-73-1	1,3-Dichlorobenzene	2.0 U
106-46-7	1,4-Dichlorobenzene	2.0 U
100-51-6	Benzyl Alcohol	2.0 U
95-50-1	1,2-Dichlorobenzene	2.5 U
95-48-7	2-Methyphenol	1.0 U
39638-32-9	bis(2-chloroisopropyl)Ether	2.5 U
106-44-5	4-Methyphenol	1.0 U
621-84-7	N-Nitroso-Di-n-Propylamine	1.5 U
67-72-1	Hexachloroethane	2.0 U
98-95-3	Nitrobenzene	2.5 U
78-59-1	Isophorone	2.5 U
88-75-5	2-Nitrophenol	2.0 U
105-67-9	2,4-Dimethyphenol	2.0 U
65-85-0	Benzoic Acid	3.0 U
111-91-1	bis(2-Chloroethoxy)Methane	2.5 U
120-83-2	2,4-Dichlorophenol	2.0 U
120-82-1	1,2,4-Trichlorobenzene	2.0 U
91-20-3	Naphthalene	2.0 U
106-47-8	4-Chloroaniline	2.0 U
87-68-3	Hexachlorobutadiene	2.5 U
59-50-7	4-Chloro-3-Methylphenol	1.5 U
91-57-6	2-Methylnaphthalene	2.0 U
77-47-4	Hexachlorocyclopentadiene	2.0 U
88-06-2	2,4,6-Trichlorophenol	1.5 U
95-95-4	2,4,5-Trichlorophenol	1.5 U
91-58-7	2-Chloronaphthalene	1.5 U
88-74-4	2-Nitroaniline	1.0 U
131-11-3	Dimethyl Phthalate	1.5 U
208-96-8	Acenaphthylene	1.5 U
99-09-2	3-Nitroaniline	2.5 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	1.5 U
51-28-5	2,4-Dinitrophenol	1.5 U
100-02-7	4-Nitrophenol	1.5 U
132-64-9	Dibenzofuran	1.0 U
121-14-2	2,4-Dinitrotoluene	1.0 U
606-20-2	2,6-Dinitrotoluene	1.0 U
84-66-2	Diethylphthalate	1.0 U
7005-72-3	4-Chlorophenyl-phenylether	1.0 U
86-73-7	Fluorene	1.0 U
100-01-6	4-Nitroaniline	3.0 U
534-52-1	4,6-Dinitro-2-Methylphenol	1.5 U
86-30-6	N-Nitrosodiphenylamine (1)	1.5 U
101-55-3	4-Bromophenyl-phenylether	1.5 U
118-74-1	Hexachlorobenzene	1.5 U
97-86-5	Pentachlorophenol	2.0 U
85-01-8	Phenanthrene	1.0 U
120-12-7	Anthracene	2.5 U
84-74-2	Di-n-Butylphthalate	2.0 U
206-44-0	Fluoranthene	1.5 U
129-00-0	Pyrene	1.5 U
85-68-7	Butylbenzylphthalate	3.5 U
91-84-1	3,3'-Dichlorobenzidine	NR
56-55-3	Benzo(a)Anthracene *	1.5 U
117-81-7	bis(2-Ethylhexyl)Phthalate	1.0 U
218-01-9	Chrysene *	—
117-84-0	Di-n-Octyl Phthalate	1.5 U
205-99-2	Benzo(b)Fluoranthene **	1.5 U
207-08-9	Benzo(k)Fluoranthene **	—
50-32-8	Benzo(a)Pyrene	2.0 U
193-39-5	Indeno(1,2,3-cd)Pyrene	3.5 U
53-70-3	Dibenz(a,h)Anthracene	2.5 U
181-24-2	Benzo(g,h,i)Perylene	4.0 U

(1)-Cannot be separated from diphenylamine

\*THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

\*\*THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

Laboratory Name S-CUBED  
Case No 9288 (3554E)

Sample Number  
LAB BLANK 4-11

Organics Analysis Data Sheet  
(Page 3)

Pesticide / PCBs

Concentration Low Medium (Circle One)  
Date Extracted / Prepared 4-11-88  
Date Analyzed 4-12-88  
Conc / Dil Factor 1.0  
Percent Moisture (decanted) NA

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug / 10 mg / Kg (Circle One)
319-84-6	Alpha-BHC	0.010 $\mu$
319-85-7	Beta-BHC	0.005 $\mu$
319-86-8	Delta-BHC	0.005 $\mu$
58-69-9	Gamma-BHC (Lindane)	0.005 $\mu$
76-44-8	Heptachlor	0.030 $\mu$
309-00-2	Aldrin	0.005 $\mu$
1024-57-3	Heptachlor Epoxyde	0.005 $\mu$
959-98-8	Endosulfan I	0.010 $\mu$
60-57-1	Dieldrin	0.010 $\mu$
72-55-9	4,4'-DDE	0.005 $\mu$
72-20-8	Endrin	0.010 $\mu$
33213-65-9	Endosulfan II	0.010 $\mu$
72-54-8	4,4'-DDD	0.020 $\mu$
1031-07-8	Endosulfan Sulfate	0.100 $\mu$
50-29-3	4,4'-DDT	0.020 $\mu$
72-43-5	Methoxychlor	0.020 $\mu$
53494-70-5	Endrin Ketone	0.030 $\mu$
57-74-9	Chlordane	0.020 $\mu$
8001-35-2	Toxaphene	0.25 $\mu$
12674-11-2	Aroclor-1016	0.10 $\mu$
11104-28-2	Aroclor-1221	NA
11141-16-5	Aroclor-1232	NA
53469-21-9	Aroclor-1242	0.10 $\mu$
12672-29-6	Aroclor-1248	0.10 $\mu$
11097-69-1	Aroclor-1254	0.10 $\mu$
11096-82-5	Aroclor-1260	0.10 $\mu$
74421-93-4	DODIN ALDENEYDE	0.030 $\mu$

$V_i$  = Volume of extract injected ( $\mu$ l)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract ( $\mu$ l)

$v_s$  1000 ml or  $w_s$  NA  $v_i$  1.00 ml  $v_t$  1.0  $\mu$ l

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Sample Number

LAB BLANK

44V2021 (VOA)

RRBAG (BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	No TIC FOUND.	VOA		
2.				
3.	No TIC Found	BNA		
4.				
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ORGANICS ANALYSIS DATA SHEET  
(Page 1)

Laboratory Name: 6-CUBED  
Lab Sample ID No: 4AV2021 (NOA), LBA02 (ABW)  
Sample Matrix: WATER  
Data Release Authorized By: [initials]

Case No: 9288 (SAS3554E)  
QC Report Nos: N.R  
Contract Nos: 6B-01-7261  
Date Sample Received: N.R

VOLATILE COMPOUNDS

Concentration: Low/Medium (Circle One)  
Date Extracted/Prepared: 04-04-88  
Date Analyzed: 04-04-88  
Conc/Dil Factor: 1.0 pH: N.R  
Percent Moisture (Not Decanted): N.R

CAS Number		ug/l or ug/kg (Circle One)	CAS Number		ug/l or ug/kg (Circle One)
74-87-3	CHLOROMETHANE	10 U	78-87-5	1,2-DICHLOROPROPANE	1.5 U
74-83-9	BROMOMETHANE	10 U	10061-02-6	TRANS-1,3-DICHLOROPROPENE	1 U
75-01-4	VINYL CHLORIDE	10 U	79-01-6	TRICHLOROETHENE	1.5 U
75-00-3	CHLOROETHANE	1.5 U	124-48-1	DIBROMOCHLOROMETHANE	1.5 U
75-09-2	METHYLENE CHLORIDE	2.2	79-00-5	1,1,2-TRICHLOROETHANE	1.5 U
67-64-1	ACETONE	75 U	71-43-2	BENZENE	1.5 U
75-15-0	CARBON DISULFIDE	3 U	10061-01-5	CIS-1,3-DICHLOROPROPENE	2 U
75-35-4	1,1-DICHLOROETHENE	1.5 U	110-75-8	2-CHLOROETHYL VINYL ETHER	1.5 U
75-34-3	1,1-DICHLOROETHANE	1.5 U	75-25-2	BROMOFORM	1.5 U
	TOTAL-1,2-DICHLOROETHENE'S	1.5 U	591-78-6	2-HEXANONE	50 U
67-66-3	CHLOROFORM	1.5 U	108-10-1	4-METHYL-2-PENTANONE	3 U
107-06-2	1,2-DICHLOROETHANE	1.5 U	127-18-4	TETRACHLOROETHENE	1.5 U
78-93-3	2-BUTANONE	50 U	79-34-5	1,1,2,2-TETRACHLOROETHANE	1.5 U
71-35-6	1,1,1-TRICHLOROETHANE	1.5 U	108-88-3	TOLUENE	1.5 U
56-23-5	CARBON TETRACHLORIDE	1.5 U	108-90-7	CHLOROBENZENE	1.5 U
108-05-4	VINYL ACETATE	15 U	100-41-4	ETHYLBENZENE	1.5 U
75-27-4	BROMODICHLOROMETHANE	1.5 U	100-42-5	STYRENE	1 U
107-02-8	ACROLEIN	100 U		M-XYLENE	2 U
107-13-1	ACRYLONITRILE	50 U		O/P-XYLENE	2.5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifier are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value: If the result is a value greater than or equal to the detection limit, report the value.
- U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (eg.10U) based on necessary concentration/dilution action. (this is not necessarily the instrument detection limit.) The footnotes should read:U=Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J Indicates an estimated value: This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (eg 10J). If limit of detection is 10 ug/L and a concentration of 3ug/L is calculated, report as 3J.

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides > or = 10 ng/ul in the final extract should be confirmed by GC/MS.

B This flag is used when analyte is found in the blank as well as sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Matrix spiked compound.

Laboratory Name S-CUBED  
Case No 9288/3554F

Sample Number  
Lab Blank 4-7

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration Low Medium (Circle One)  
Date Extracted/Prepared 4-7-88  
Date Analyzed 4-10-88  
Conc 'Dil Factor 1.0  
Percent Moisture (Decanted) N.A.

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	2.0 U
111-44-4	bis(2-Chloroethyl)Ether	1.5 U
95-57-8	2-Chlorophenol	2.0 U
541-73-1	1, 3-Dichlorobenzene	2.0 U
106-46-7	1, 4-Dichlorobenzene	2.0 U
100-51-6	Benzyl Alcohol	2.0 U
85-50-1	1, 2-Dichlorobenzene	2.5 U
95-48-7	2-Methylphenol	1.0 U
39638-32-9	bis(2-chloroethyl)Ether	2.5 U
106-44-5	4-Methylphenol	1.0 U
621-64-7	N-Nitroso-Di-n-Propylamine	1.5 U
67-72-1	Hexachloroethane	2.0 U
98-95-3	Nitrobenzene	2.5 U
78-59-1	Iso phorone	2.5 U
88-75-5	2-Nitrophenol	2.0 U
105-67-9	2, 4-Dimethylphenol	2.0 U
65-85-0	Benzoic Acid	3.0 U
111-91-1	bis(2-Chloroethyl)Methane	2.5 U
120-83-2	2, 4-Dichlorophenol	2.0 U
120-82-1	1, 2, 4-Trichlorobenzene	2.0 U
91-20-3	Naphthalene	2.0 U
106-47-8	4-Chloroaniline	2.0 U
87-68-3	Hexachlorobutadiene	2.5 U
59-50-7	4-Chloro-3-Methylphenol	1.5 U
91-57-6	2-Methylnaphthalene	2.0 U
77-47-4	Hexachlorocyclopentadiene	2.0 U
88-08-2	2, 4, 6-Trichlorophenol	1.5 U
95-95-4	2, 4, 5-Trichlorophenol	1.5 U
91-58-7	2-Chloronaphthalene	1.5 U
88-74-4	2-Nitroaniline	1.0 U
131-11-3	Dimethyl Phthalate	1.5 U
208-96-8	Acenaphthylene	1.5 U
89-09-2	3-Nitroaniline	2.5 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	1.5 U
51-28-5	2, 4-Dinitrophenol	1.5 U
100-02-7	4-Nitrophenol	1.5 U
132-64-9	Dibenzofuran	1.0 U
121-14-2	2, 4-Dinitrotoluene	1.0 U
606-20-2	2, 6-Dinitrotoluene	1.0 U
84-86-2	Diethylphthalate	2.1
7005-72-3	4-Chlorophenyl-phenylether	1.0 U
88-73-7	Fluorene	1.0 U
100-01-8	4-Nitroaniline	3.0 U
534-52-1	4, 6-Dinitro-2-Methylphenol	1.5 U
86-30-6	N-Nitrosodiphenylamine (1)	1.5 U
101-55-3	4-Bromophenyl-phenylether	1.5 U
118-74-1	Hexachlorobenzene	1.5 U
87-86-5	Pentachlorophenol	2.0 U
85-01-8	Phenanthrene	1.0 U
120-12-7	Anthracene	2.5 U
84-74-2	Di-n-Butylphthalate	2.0 U
206-44-0	Fluoranthene	1.5 U
129-00-0	Pyrene	1.5 U
85-88-7	Butylbenzylphthalate	3.5 U
91-84-1	3, 3'-Dichlorobenzidine	NR
56-55-3	Benz(a)Anthracene *	1.5 U
117-81-7	bis(2-Ethyhexyl)Phthalate	1.0 U
218-01-9	Chrysene *	—
117-84-0	Di-n-Octyl Phthalate	1.5 U
205-99-2	Benz(b)Fluoranthene # *	1.5 U
207-08-9	Benz(k)Fluoranthene # *	—
50-32-8	Benz(a)Pyrene	2.0 U
193-39-5	Indeno[1, 2, 3-cd]Pyrene	3.5 U
53-70-3	Dibenzo[ <i>a, h</i> ]Anthracene	2.5 U
191-24-2	Benz[ <i>a, h, i</i> ]Perylene	4.0 U

(1)-Cannot be separated from diphenylamine

\* THESE TWO PARAMETERS ARE REPORTED AS A TOTAL

# THESE TWO PARAMETERS ARE REPORTED AS A TOTAL  
Form I

Laboratory Name S-CUBED  
Case No 9288 (SAS3554E)

Sample Number

LAB BLANK

4AV2021(VOA)

LB A07 (BNA)

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l) or ug/kg)
1.	No TIC FOUND.	VOA		
2.				
3. 21964498	1,13 tetradecadiene	BNA	1294	22 J
4. 19812647	1,14 tetradecane diol	BNA	1298	6 J
5.				
6.				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
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20.				
21.				
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23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				



# ORGANIC TRAFFIC REPORT

(FOR CLP USE ONLY)

TYPE OF ACTIVITY (CIRCLE ONE) <input checked="" type="radio"/> SUPERFUND <input type="radio"/> OTHER <input type="radio"/> OAM OTHER <input type="radio"/> DMR		SHIP TO: <b>5-Cubed (53)</b>	DATE REC'D: <b>5/29/88</b>	SDG NO: <b>851</b>
NON-SUPERFUND <input type="radio"/> PROGRAM		3398 La Sierra Mountain Rd San Diego, CA 92121 ATTN: <u>Elaine Walters</u>	REC'D BY: <u>Elaine Walters</u>	
SAMPLE DESCRIPTION (ENTER IN BOX ORANGE, GREEN, OR BLUE) 1. SURFACE WATER <input type="checkbox"/> 2. GROUND WATER <input type="checkbox"/> 3. LEACHATE <input type="checkbox"/> 4. SOIL <input type="checkbox"/> 5. WASTE (TRASH) <input type="checkbox"/> 6. OIL (BAS) <input type="checkbox"/> 7. GASES (TRUCKS 2V2) <input type="checkbox"/> 8. UNKNOWN <input type="checkbox"/>		SAMPLING DATE: <b>BEGIN: 3/29/88 END: 3/31/88</b>	LABORATORY CONTRACT NO. <b>7261</b>	UNIT PRICE
REGION NO: <b>T</b>	SAMPLING COMPANY <b>MPCA</b>	DATE SHIPPED: <b>4/1/88</b>	RECEIVED BY:	DATE REC'D:
SAMPLER (NAME) <b>REBECCA LOFGREN</b>		AIRBILL NO: <b>6820695814</b>	CONTRACT NO./PRICE:	

SAMPLE ID	NUMBER (FROM LABELS)	SAMPLE BOX		SAMPLE CONDITION	HIGH CON PHASES	SOLID	WATER	LIQUID	NON WATER
		CONT.	CONT.						
ET 831	2	X	X	H-4					
ET 832	2	L	X X X	H-3					
ET 833	2	X	X X X	H-1					
ET 834	2	L	X X X	H-2					
ET 835	2	X	X X X	H-1					
ET 836	2	L	X X X	H-7					
ET 837	2	L	X X X	H-5					
ET 838	2	L	X X X	H-8					
ET 839	2	L	X X X	H-9					
ET 840	2	L	X	TRIP BANK					

SAMPLING COMPLETE  
(1 of 4 coolers shipped)

IDENTIFICATION  
OF NAMES OF INTERNAL STANDARD (IS) AND  
SURROGATES (SURR) ON RIC SHEET

VOLATILE

IS1 : BROMOCHLOROMETHANE  
IS2 : 1,4-DIFLUOROBENZENE  
IS3 : D5-CHLOROBENZENE  
SURR1 : D8-TOLUENE  
SURR2 : D4-1,2-DICHLOROETHANE  
SURR3 : BROMOFLUOROBENZENE

SEMI-VOLATILE

IS1 : D4-DICHLOROBENZENE  
IS2 : D8-NAPHTHALENE  
IS3 : D10-ACENAPHTHENE  
IS4 : D10-PHENANTHRENE  
IS5 : D12-CHRYSENE  
IS6 : D12-PERYLENE  
SURR3 : D5-NITROBENZENE  
SURR2 : D5-PHENOL  
SURR1 : 2-FLUOROPHENOL  
SURR4 : 2-FLUOROBIPHENYL  
SURR5 : 2,4,6-TRIBROMOPHENOL  
SURR6 : D14-TERPHENYL

## GC/MS TUNING AND MASS CALIBRATION

## **Decafluorotriphenylphosphine (DFTPP)**

Case No. 9288/3554E Contractor S-Cubed Contract No. 6801-72(a).

Instrument ID HP1. Date / Time 4/10/88 1:50

Lab ID :TA07C:02 Data Release Authorized By: \_\_\_\_\_ b)

a/z	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	42.77 OK
68	less than 2.0% of mass 69	0.00 OK ( 0.00) \$1
69	mass 69 relative abundance	56.64
70	less than 2.0% of mass 69	0.00 OK ( 0.00) \$1
127	40.0 - 60.0% of mass 198	47.44 OK
197	less than 1.0% of mass 198	.93 OK
198	base peak, 100% relative abundance	100.00 OK
199	5.0 - 9.0% of mass 198	6.76 OK
225	10.0 - 30.0% of mass 198	20.12 OK
365	greater than 1.00% of mass 198	2.06 OK
441	present, but less than mass 443	9.44 OK
442	greater than 40.0% of mass 198	62.96 OK
443	17.0 - 23.0% of mass 442	13.46 OK (21.37) \$2

**THIS PERFORMANCE TUNE APPLIES TO THE  
FOLLOWING SAMPLES, BLANKS AND STANDARDS.**

#1 - Value in parenthesis is % mass 69.

#2 - Value in parenthesis is % mass 442.

## GC/MS TUNING AND MASS CALIBRATION

## Decafluorotriphenylphosphine (DFTPP)

Case No. 9293/3554E - Contractor S - CUBED Contract No. 6801-7261

Instrument ID #1(HP) Date / Time 4/13/88 17:27

Lab ID >TM13A::D3 Data Release Authorized By: [Signature]

m/z	ION ABUNDANCE CRITERIA	RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	40.93 OK
68	less than 2.0% of mass 69	0.60 OK ( 0.00 ) #1
69	mass 69 relative abundance	55.13
70	less than 2.0% of mass 69	0.00 OK ( 0.00 ) #1
127	40.0 - 60.0% of mass 198	46.04 OK
197	less than 1.0% of mass 198	0.00 OK
198	base peak, 100% relative abundance	100.00 OK
199	5.0 - 9.0% of mass 198	7.42 OK
275	10.0 - 30.0% of mass 198	21.56 OK
365	greater than 1.00% of mass 198	2.10 OK
441	present, but less than mass 443	10.14 OK
442	greater than 40.0% of mass 198	64.91 OK
443	17.0 - 23.0% of mass 442	14.67 OK (22.59) #2

THIS PERFORMANCE TUNE APPLIES TO THE  
FOLLOWING SAMPLES, BLANKS AND STANDARDS.#1 - Value in parenthesis is % mass 69.  
#2 - Value in parenthesis is % mass 442.

SAMPLE ID	LAB_ID	DATE_OF_ANALYSIS	TIME_OF_ANALYSIS
20 ppm TGA1	SF 301	4-13-88	1753 ✓
50 ppm TGA1	SF 302		1853 ✓
100 ppm TGA1	SF 303		1951 ✓
Lab Blk 4-4	RRBAG		2051 ✓
ET 726	RR 726		2150
E 722	RR 722		2249
EW 274	RR 274	4-13-88	2348
EW 275	RR 275	4-14-88	0047
EW 280	RR 280		0146
ET 837	ET 837		0244 ✓
ET 838	ET 838		0342 ✓

## GC/MS TUNING AND MASS CALIBRATION

## **Decafluorotriphenylphosphine (DFTPP)**

Case No. 928.8 / 3554E

Contractor S - CUBED

Contract No. 6801-7261

Instrument ID #1(HP)

Date / Time 4/14/88 15:56

Lab ID >TA14A::02

**Date Release Authorized By:**

M/Z	ION ABUNDANCE CRITERIA	RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	36.79 OK
68	less than 2.0% of mass 69	0.00 OK ( 0.00) \$1
69	mass 69 relative abundance	48.04
70	less than 2.0% of mass 69	0.00 OK ( 0.00) \$1
127	40.0 - 60.0% of mass 198	41.97 OK
197	less than 1.0% of mass 198	0.00 OK
198	base peak, 100% relative abundance	100.00 OK
199	5.0 - 9.0% of mass 198	7.33 OK
275	10.0 - 30.0% of mass 198	20.61 OK
365	greater than 1.00% of mass 198	1.77 OK
441	present, but less than mass 443	12.64 OK
442	greater than 40.0% of mass 198	80.91 OK
443	17.0 - 23.0% of mass 442	18.33 OK (22.66) \$2

**THIS PERFORMANCE TUNE APPLIES TO THE  
FOLLOWING SAMPLES, BLANKS AND STANDARDS.**

§1 - Value in parenthesis is % mass 69.

#2 - Value in parenthesis is % less 442.

## GC/MS TUNING AND MASS CALIBRATION

#### **Decafluorotriphenylphosphine (DFTPP)**

Case No. 9288/3554E Contractor S - CUBED Contract No. 6801-7261

Instrument ID #1(HP) Date / Time 4/19/88 10:02

Lab ID :TA19A:D2 Data Release Authorized By: WJ

m/z	ION ABUNDANCE CRITERIA		%RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198		51.50 OK
68	less than 2.0% of mass 69		1.16 OK (1.003) #1
69	mass 69 relative abundance		64.09
70	less than 2.0% of mass 69		0.00 OK (0.00) #1
127	40.0 - 60.0% of mass 198		51.09 OK
197	less than 1.0% of mass 198		0.00 OK
198	base peak, 100% relative abundance		100.00 OK
199	5.0 - 9.0% of mass 198		6.91 OK
275	10.0 - 30.0% of mass 198		18.55 OK
365	greater than 1.00% of mass 198		2.09 OK
441	present, but less than mass 443		7.13 OK
442	greater than 40.0% of mass 198		45.09 OK
443	17.0 - 23.0% of mass 442		9.12 OK (20.24) #2

**THIS PERFORMANCE TUNE APPLIES TO THE  
FOLLOWING SAMPLES, BLANKS AND STANDARDS.**

**81 - Value in parenthesis is % pass 69.**

f2 - Value in parenthesis is % mass 442.

## GC/MS TUNING AND MASS CALIBRATION

## Bromo-fluorobenzene (BFB)

Case No. 92K 9288 Contractor S-CUBED Contract No 58-01-5862  
Instrument ID 1G2 Date 20 Mar 00 ✓ Time 0159 ✓  
Lab ID JZBNZBFB ✓ Data Release Authorized By: MJ

ratio	ION ABUNDANCE CRITERIA	RELATIVE ABUNDANCE
50	19.0 - 40.0% of the base peak	21.9 ✓
73	30.0 - 50.0% of the base peak	53.5 ✓
93	Base peak, 100% relative abundance	100 ✓
96	5.0 - 9.0% of the base peak	6.90 ✓
173	Less than 1.0% of the base peak	no peak ✓
174	Greater than 90.0% of the base peak	57.8 ✓
175	5.0 - 9.0% of mass 174	4.07 ✓ (7.04) ✓
176	Greater than 95.0%, but less than 101.0% of mass 174	56.5 ✓ (97.8) ✓
177	5.0 - 9.0% of mass 176	3.80 ✓ (4.73) ✓

**THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING  
SAMPLES, BLANKS AND STANDARDS.**

<sup>1</sup>Value in parentheses is % man 174

<sup>2</sup>Yield in percentage is 3 times 178.

## **GC/MS TUNING AND MASS CALIBRATION**

## Bromofluorobenzene (BFB)

Case No. 9288 Contractor S-CUBED Contract No 58-31-5862  
Instrument ID VG 2 Date 4-APR-88 ✓ Time 19:46 ✓  
Lab ID A4V2BFB ✓ Data Release Authorized By: pn

mass	ION ABUNDANCE CRITERIA	RELATIVE ABUNDANCE
50	15.0 - 42.0% of the base peak	19.9 ✓
73	32.0 - 62.0% of the base peak	59.5 ↙
95	Base peak, 100% relative abundance	100 ↘
96	5.0 - 9.0% of the base peak	7.06 ↙
173	Less than 1.0% of the base peak	none
174	Greater than 50.0% of the base peak	56.7 ✓
175	5.0 - 9.0% of mass 174	3.89 ✓ (6.8) <sup>1</sup>
176	Greater than 25.0%, but less than 101.0% of mass 174	54.9 ✓ (96.8) <sup>1</sup>
177	5.0 - 9.0% of mass 176	3.66 ✓ (6.56) <sup>2</sup>

**THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING  
SAMPLES, BLANKS AND STANDARDS.**

<sup>1</sup>Value in parentheses is % mass 174

<sup>2</sup>Value in parenthesis is % mass 175.

## **GC/MS TUNING AND MASS CALIBRATION**

## **Bromo-*o*-fluorobenzene (BFB)**

Gas No. 1288

10

S-CUBED

## Contract

58-31-9286

Instrument 10 VG2

100

5-1985

10

— 1 —

4AV2BFB

• 100

10 of 10

#### **NEW AIRLINE AND CARRIER CRITERIA**

### **RELATIVE ABUNDANCE**

50	15.0 - 40.0% of the base peak	10.0 ✓
73	30.0 - 60.0% of the base peak	52.5 ✓
95	Base peak, 100% relative abundance	100 ✓
96	5.0 - 9.0% of the base peak	6.8D ✓
173	Less than 1.0% of the base peak	No peak ✓
174	Greater than 50.0% of the base peak	59.3 ✓
175	5.0 - 9.0% of mass 174	4.16 ✓ (7.0)'
176	Greater than 95.0%, but less than 101.0% of mass 174	58.7 ✓ (99.0)'
177	5.0 - 9.0% of mass 176	37+ 3.65 fm (6.385 <sup>2</sup> ) fm 6.333 fm

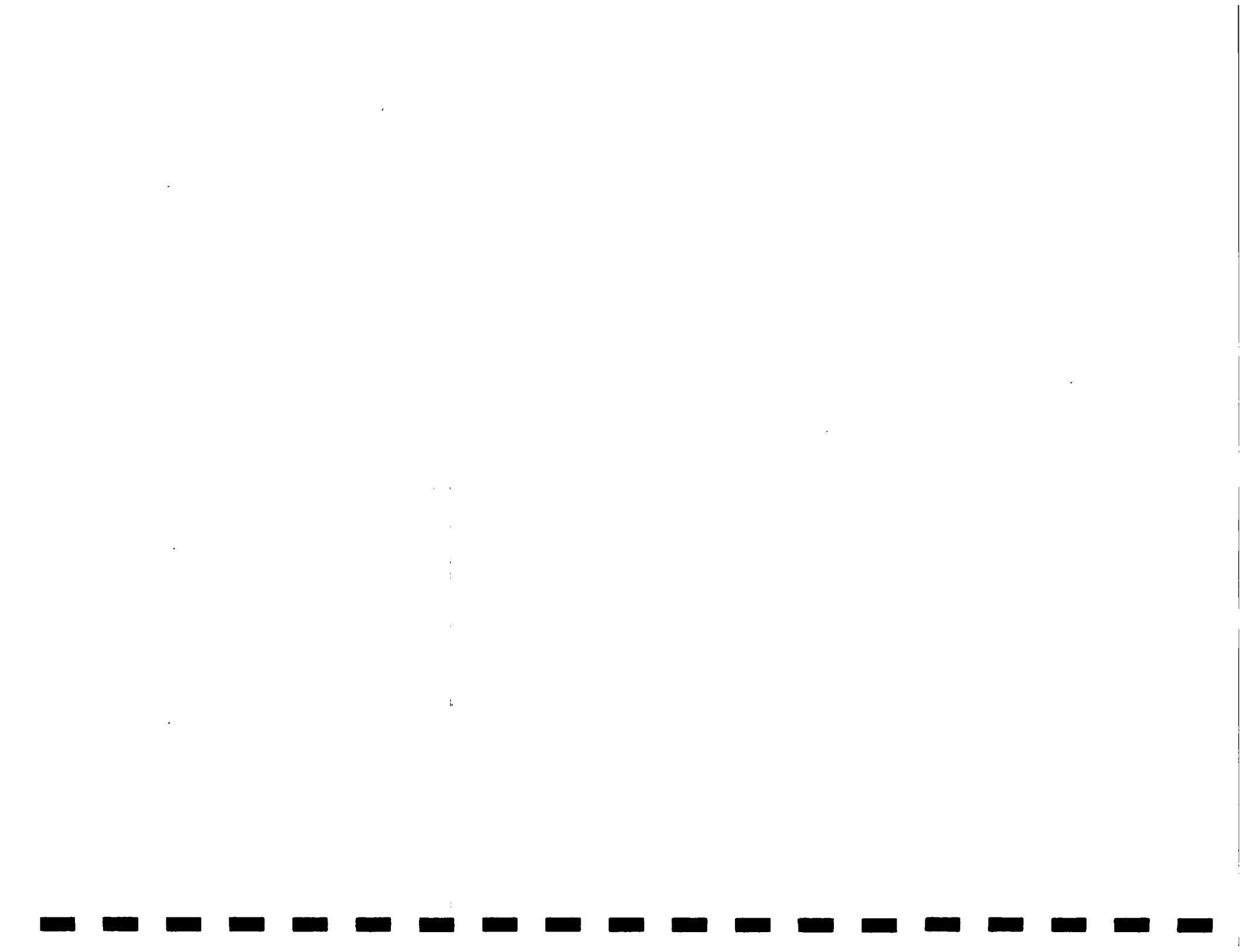
**THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING  
SAMPLES, BLANKS AND STANDARDS.**

<sup>1</sup>Value in parenthesis is % mass 174  
<sup>2</sup>Value in parenthesis is % mass 175.

<sup>2</sup> Values in parentheses are mass 176.

# **INORGANIC DATA**

# **INORGANIC DATA**



SF 5011  
MPCA

## COVER SHEET

(MT)

LABORATORY RESPONSE TO RESULTS OF  
CONTRACT COMPLIANCE SCREENING (CCS)~~Under G.W.~~

Response To: (check one)

Organics CCS

Inorganics CCS

Response materials sent to Organics CCS should be sent to the attention of Doris Ling, SMO.

Response materials sent to Inorganics CCS should be sent to the attention of Sa'ad Masri, SMO.

Laboratory Name: Rocky Mountain Analytical

Response Date 5-18-88EPA Contract No: 68-01-743Ko

SDG No:

MEW615Date Screening  
Results Received  
at Laboratory

Case No:

92885-10-88

Sample Nos.\*

SAS 3554E


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\* Only list sample numbers that require reconciliation.

This form is used to identify materials sent in response to results of Contract Compliance Screening (CCS). A separate form must accompany the response for each Case.

Please indicate (on the attached continuation form) which fractions and/or which criteria correspond with your resubmission. Respond materials sent to CCS should also be copied to the Region and to EMSL/LV, each with this blue Cover Sheet.

# INORGANICS

## LABORATORY RESPONSE TO RESULTS OF CCS

Criterion	Comments
B-H	Data packages are to be screened for RAS requirements only.
T	Submitted traffic report for this case.
FB2	Resubmitted Form IVB with "NR" for Sb, Cd, and Zn.
J	Refer to item #6 in attached addendum. Also note that Form IX is for ICP serial dilutions.
J2	Serial dilution results were done at a 5x dilution; resubmitted raw data with "5x" entered.
R	Refer to item #15 in attached addendum.
I	Data packages are to be screened for RAS requirements only.
R2	See above.

**ADDENDUM TO  
LABORATORY RESPONSE TO RESULTS OF CCS**

1. The contract states that values greater than 10 are to be reported to 3 significant figures. Raw data often have more significant figures.  
(See the January 25, 1988 letter from Tony Maiorana to the USEPA's Project Officer, p.2, item 3.)
2. This is not required or even mentioned in the contract. In fact, this is impossible when lead is analyzed by both ICP and AA on different samples in the same case.  
(See p.1, item 2 in the above mentioned letter.)
3. We submit CCVs only for that case reported. Earlier CCVs are not required in the data package.  
(See p.3, item 17 in the above mentioned letter.)
4. There are no Calcium correction factors on this ICP.  
(See p.2, item 12 in the above mentioned letter.)
5. Cyanide is being prepared using a micro-distillation process, according to the historical analysis instituted in 1981. Both the Project Officer and the Contract Officer are studying this point and will make a determination at some future time.
6. A blank field under the M column is considered an entry and indicates that the serial dilution was not required because the analyte was not analyzed by ICP.  
(SOW787, B-29)
7. Beginning 3-10-88 ICV/ICB and CCV/CCB times for Mercury are reported.
8. The % solids of the sample is used to calculate the duplicate value. (See p.D-84 of SOW787, paragraph 5, last sentence.)
9. Enseco has recognized a problem with the furnace analysis of the solid LCS sample. The contract does not allow an MSA determination for the solid or aqueous LCS result (p. E-14). This causes problems for the correct quantitation of the solid LCS because the true values and control limits were set using results determined under SOW784 and SOW785 which do allow MSA quantitation of the LSC result. Enseco has contacted our Project Officer, Bill Langley, concerning this issue. Until further notification, he has instructed us to perform MSA analyses on the furnace solid LCS following the sample analysis decision tree in SOW787.
10. As required by the contract, the reported concentrations in the ICP raw data are instrument output from the diluted sample.

ADDENDUM  
Page Two

11. The dilution factor is reported for all samples and ICV's. Standard dilution factor for CCV's is referenced at the beginning of data package on standard sources page. Also, the dilution factor is being misread by CCS screeners as volume.
12. Some of the modifications requested in the 12/87 letter of contract revisions from the Contract Officer are being instituted while others are being reconsidered by the Contract Officer. At this time, there has not been a final determination by the Contract Officer to institute any of these changes.
13. There is no contract reference given. In order to respond, a specific contract reference must be provided.
14. As per the April 7, 1988 letter from Contract Officer Mary Stotler to Enseco, CCS was found to be in error and have been directed to remove this item from the screens.
15. The question of significant figures is currently being discussed by the Contract Officer and Project Officer; as of yet the issue has not been resolved. Please also note that the values on Forms II and III are taken directly from instrument readout. There is a hidden digit which does not print out on furnace tapes, therefore it appears that the rounding is being done incorrectly.



USEPA CONTRACT LABORATORY PROGRAM  
SAMPLE MANAGEMENT OFFICE  
P.O. BOX 818 ALEXANDRIA, VA 22313  
703/557-2490 FTS-557-2490

CASE NO: 9288

SAS NO:  
(IF APPLICABLE) 3554-E

## INORGANIC TRAFFIC REPORT

(FOR CLP USE ONLY)

TYPE OF ACTIVITY (CIRCLE ONE)		SHIP TO:	DATE REC'D:	SDG NO:
SUPERFUND—PA SI ES RIFS RD RA ER • NPL & O&M OTHER		ENSECO/Rocky Mountain Area 4955 Willow Street Arvada, CO 80002 ATTN: Kathy Hutchinson	REC'D BY: <i>Linda Dellaire</i>	442-88- MEW 615 IFB+250
NON-SUPERFUND— PROGRAM		SAMPLING DATE: BEGIN: 3/29/88 END: 3/31/88	LABORATORY CONTRACT NO.	UNIT PRICE
SAMPLE DESCRIPTION (ENTER IN BOX A) 1. SURFACE WATER 2. GROUND WATER 3. LEACHATE 4. SOIL 5. SEDIMENT 6. OIL (SAS)		DATE SHIPPED: 4/1/88 CARRIER: Fed.	68-01706	
REGION NO: 10 SAMPLING COMPANY MPCA		AIRBILL NO: 6820695851	REC'D BY:	CONTRACT NO./PRICE:
SAMPLER: (NAME) REBECCA LOFGREN				

SAMPLE NUMBER (FROM LABELS)	DESCRIPTION (FROM BOX 1)	CONCENTRATION LOW, M, MED, H	RAS ANALYSIS				SPECIAL HANDLING	STATION LOCATION	SAMPLE CONDITION ON RECEIPT	HIGH CONC. PHASES (CHECK)			
			TOTAL METALS	PCVANIDE	DISSOLVED METALS	HIGH PH. ONLY (SAS)							
											SOLID	WATER-MIS LIQUID	NON WATER - MIS LIQUID
MEW 615	2 L XX							H-4	OK				
MEW 616	2 L XX							H-3					
MEW 617	2 L XY							H-1					
MEW 618	2 L XX							H-2					
MEW 619	2 L XX							H-6					
MEW 620	2 L XX							H-7					
MEW 621	2 L XX							H-5					
MEW 622	2 L XY							H-8					
MEW 623	2 L XY <i>SDS-Final Sample</i>							H-9					

SAMPLING COMPLETE

1 of 2 coolers Shipped

5B  
POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476mew616ALab Code: ENSECOCase No.: 9288SAS No.: 35546SDG No.: mew615Matrix (soil/water): WATERLevel (low/med): LOW

Concentration Units: ug/L

Analyte	Control Limit %R	Spiked Sample Result (SSR) C	Sample Result (SR) C	Spike Added (SA)	%R	Q	M
Aluminum						NR	
Antimony						NP	
Arsenic						NR	
Barium						NR	
Beryllium						NR	
Cadmium						NR	
Calcium						NR	
Chromium		19.3	4.0	20.0	96.5	NR	
Cobalt						NR	
Copper						NR	
Iron						NR	
Lead						NR	
Magnesium						NR	
Manganese						NR	
Mercury						NR	
Nickel						NR	
Potassium						NR	
Selenium						NR	
Silver						NR	
Sodium						NR	
Thallium						NR	
Vanadium						NR	
Zinc						ZNR	
Cyanide						NR	

Comments:

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卷之三

	CE	CF
1	1.000000	1.000000
2	1.000000	1.000000
3	1.000000	1.000000
4	1.000000	1.000000
5	1.000000	1.000000
6	1.000000	1.000000
7	1.000000	1.000000
8	1.000000	1.000000
9	1.000000	1.000000
10	1.000000	1.000000
11	1.000000	1.000000
12	1.000000	1.000000
13	1.000000	1.000000
14	1.000000	1.000000
15	1.000000	1.000000
16	1.000000	1.000000
17	1.000000	1.000000
18	1.000000	1.000000
19	1.000000	1.000000
20	1.000000	1.000000
21	1.000000	1.000000
22	1.000000	1.000000
23	1.000000	1.000000
24	1.000000	1.000000
25	1.000000	1.000000
26	1.000000	1.000000
27	1.000000	1.000000
28	1.000000	1.000000
29	1.000000	1.000000
30	1.000000	1.000000
31	1.000000	1.000000
32	1.000000	1.000000
33	1.000000	1.000000
34	1.000000	1.000000
35	1.000000	1.000000
36	1.000000	1.000000
37	1.000000	1.000000
38	1.000000	1.000000
39	1.000000	1.000000
40	1.000000	1.000000
41	1.000000	1.000000
42	1.000000	1.000000
43	1.000000	1.000000
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88	1.000000	1.000000
89	1.000000	1.000000
90	1.000000	1.000000
91	1.000000	1.000000
92	1.000000	1.000000
93	1.000000	1.000000
94	1.000000	1.000000
95	1.000000	1.000000
96	1.000000	1.000000
97	1.000000	1.000000
98	1.000000	1.000000
99	1.000000	1.000000
100	1.000000	1.000000

	SE	S105	AG	NA	SR	T1	SN	V	ZN
EXP# 1	-1.125	16.81	-0.0140	-0.2670	.1442	-0.0120	-0.0152	-0.0027	-0.4856
EXP# 2	-2.316	17.75	-0.0140	5.552	.0562	-0.0012	-0.0493	-0.0007	-0.4270
EXP# 3	-1.6446	17.08	-0.0140	2.677	.1017	-0.0065	-0.0224	-0.0017	-0.4862
EXP# 4	-1.6446	17.08	-0.0140	2.677	.1017	-0.0065	-0.0224	-0.0017	-0.4862

ACT NAME : ESTATE TIME : 12-4-FF-38 SAMPLE NAME: COVZ

	I-5	AL	SB	AS	BA	BE	E	CD	CA	DR
EXP# 1	1000.	.3867	.155	1.008	.9554	.9512	.9528	.957	104.7	.9291
EXP# 2	1000.	1.000	.156	1.007	1.000	1.001	1.025	.9522	104.9	.9306
AVERAGE	1000.	.9935	.155	1.007	1.001	.9932	.9525	.9527	104.7	.9295

# Enseco

January 25, 1988

William Langley  
USEPA  
401 M Street, S.W.  
Washington, D.C. 20460

Dear Bill:

Enseco has been getting 0% recommendations for payment recently on data packages submitted by Rocky Mountain Analytical Laboratory under SOW 787. In our recent telephone conversation concerning this issue you informed me that you have been led to believe that there are serious technical deficiencies with these data packages. I repeat what I told you then. The majority of the "real" problems associated with these data packages are minor and involve nothing more than report writing glitches in the software. Debugging of the software is an ongoing process that began with the generation of the first data package. This process will be completed in the near future.

The following addresses one of our recent screens on Case 8431 on a point by point basis. Most of our other screens have the same or similar comments. There are no major problems, certainly none requiring 0% payment. Many of the things screened are either not in the contract or are wrong and are indicated with an asterisk (18 out of 28 comments). The remainder would have been caught by the laboratory during debugging if a SOP had been received containing detailed acceptance criteria for the contract.

Case 8431.

- 1) CCS: "on Form I; Ca, Cu, on MDG799 don't need 8 under C column; and Ba on MDG 800."

Response: Form I's resubmitted without "8" for Ca and Cu on MDG799 and for Ba on MDG 800.

- 2) CCS: "CA1. No need to separate furnace from IC? values for IC?/CC? on Form IIA."

Response: This is not required or even mentioned in the contract. In fact, this is impossible when lead is analyzed by both IC? and AA on different samples in the same case.

William Langley  
January 25, 1988  
Page Two

- \*3) CCS: "C2. The ICY and CCY values for ICP don't agree with the raw data for the significance figures."  
Response: The contract states report to 3 significant figures for numbers greater than 10. Raw data often have more significant figures.
- \*4) CCS: "C31: missing CRDL value and %R for As and Tl on form II3."  
Response: The CRDL standards were not required since one of the standards used in the calibration was at the CRDL.
- \*5) CCS: "CB2: the initial CRDL value for Sb should be 151.9 not 152 for proper significant figures."  
Response: Same as 3).
- \*6) CCS: "D1. no need to separate the furnace and ICP values for ICS/CCS."  
Response: Same as 2)
- \*7) CCS: "E1. on form IV; the initial ICS and final ICS values for sol(AB) don't agree with the raw data (Report to one decimal place)."  
Response: Same as 3).
- 8) CCS: "F1. use an "s" for EPA sample #; (MDQ 530S)."  
Response: We are now adding "S" to the spike sample number.
- 9) CCS: "I2. Missing "+" flag for Pb on MDG5300 on form VIII."  
Response: Resubmitted Form VIII with "+" flag for Pb on sample MDG5300.
- 10) CCS: "I3. please change the EPA sample no MDQ 530 to MDG530 and MCQ 5300 to MDG5300."  
Response: Resubmitted Form VIII with corrected sample numbers.
- 11) CCS: "J1. Use an "L" for EPA Sample #."  
Response: We are now adding a "L" for the serial dilution QC sample.
- \*12) CCS: "M1. missing correction values factors for Ca.  
Response: There are no Ca correction factors required on this ICP.
- 13) CCS: "R1. the numbering of a case package must be sequential."  
Response: We have always sequentially numbered the data package and then the raw data again starting from 001. We are now numbering the entire package sequentially.

William Langley  
January 25, 1988  
Page Three

- \*14) CCS: "R2. All furnaces LCS; your raw data shows your values under PCB (ug/L) not mg/kg; final concentration shows it's mg/kg but same value under ppb is reported under final conc if these values are ug/L as in raw data your LCS is going to be out of control and samples are non-compliant; please explain and resubmit correct values with correct units in raw data and in Form VII."
- Response: The values are reported in mg/kg in the furnace raw data unless specifically stated otherwise. The values are correct as is.
- \*15) CCS: "R3. for all furnaces; under D/F (dilution factor) report one if no dilution is needed; and don't put the volumes under the D/F."
- Response: The dilution factor is reported for all samples and CCV's. Standard dilution factor for CCV's is referenced at the beginning of data package on standard sources page. Also, the dilution factor is being misread by CCS screeners as volume.
- \*16) CCS: "T1. missing SDG cover page for traffic report."
- Response: The SDG cover sheets are sent to SMO with the traffic reports. There is no contract requirement to also include them in the data package.
- \*17) CCS: "C13. need to run a CCB/CCV at end of CRI and ICSA, ICSA8 at final."
- Response: We submit CCV's only for that case reported. Earlier CCV's for other cases are not required in the data package.
- 18) CCS: "R4. in ICP raw data page 16 is out of order correct paginating your page; this page must be at the end of ICP raw data. No Action".
- Response: We are now putting this page at the end of ICP raw data.
- 19) CCS: "E2. Remove the True value for T1 Sol(A8)."
- Response: Resubmitted Form IV without the true value for T1.
- \*20) CCS: "C14. The time between ICY (8.06) and CCY (11.20) exceeds the 2 hr. limit, therefore the ICS is non-compliant."
- Response: Same as 17). CCV's are run constantly for other cases but only CCV's surrounding that case is submitted in the data.
- \*21) CCS: "F12. on form VA (SSR) for ICP don't agree with the raw data for the proper significance level."
- Response: Same as 3).
- \*22) CCS: "G1. State control limits for values < 5X CRDL. on Form VI."
- Response: No control limits are required if both values are below CRDL.

William Langley  
January 25, 1988  
Page Four

- 23) CCS: "G2. Correct control limits for % solid on Form VI."  
Response: We resubmitted Form VI with correct control limits for % solids. This was a report writer glitch which we have now fixed.
- \*24) CCS: "G3. the duplicate values on Form VI; for ICP don't agree with the raw data."  
Response: Same as 3).
- 25) CCS: "G3. need "D" for EPA sample number (MDG5300) on Form VI."  
Response: We are now adding "D" to sample number for duplicates.
- \*26) CCS: "H1. No need to separate the furnace and ICP values for LCS."  
Response: Same as 2).
- \*27) CCS: "H2. the LCS values for ICP don't agree with the raw data."  
Response: Same as 3).
- \*28) CCS: "I1. You should have value for As on MDG56300 and Pb on MDG5300 (no 0.00)."  
Response: We were not reporting values if they were not used for the concentration determination. We are now reporting values for the final concentration on Form VIII even when the value is not used.

As you can see, most of the "problems" are very minor. All of the things that were wrong (small details with no effect on the data) we have corrected. Some of the things (ie., separating ICP and furnace data) were discussed and approved with Saad Masri in Las Vegas in order to set up the Enseco report writer.

This entire situation has been exacerbated by the fact that Enseco under direct pressure from you accepted our contract minimums immediately upon award of the untested SOW 787 contract. Enseco also accepted samples beyond our minimums after persuasion and pressure was exerted by SMO.

Obviously, CCS has changed from the process set-up originally by the EPA. At least two Regions have called us and stated that CCS seems to be getting carried away with Enseco's data packages because the data packages look good to them. They did not expect perfect data packages initially under SOW 787, but are confident the deficiencies will be worked out in short order, just as in the past. To consider a data package grossly non-compliant because of minute details hurts the CLP because the compliant data percentage goes down affecting the CLP reputation, it hurts the Regional cases because it makes the data appear bad, and it hurts us because we pride ourselves on providing quality data.

William Langley  
January 25, 1988  
Page Five

Enseco is willing to send people to meet with you to address the technical problems with our data package but at this point we do not believe there are any significant major problems.

Enseco cannot indefinitely continue to analyze samples while 0% payment is being recommended. I believe you will agree that after reviewing this screen, Enseco's data packages are providing quality data that are substantially compliant with the contract and we hope you will make a fair determination of our services. At a minimum, I recommend 100% provisional payment until GCS has worked out the problems in their operation and the payment recommendations truly reflect the value of the data package to the government.

Sincerely,

*Tony*

Anthony P. Maiorana  
Enseco, ESM

cc: Mike Carter, USEPA  
Mary Stotler, USEPA  
Peter Isaacson, SMO  
Dick Thacker, SMO  
Don Trees, SMO  
Mark Carter, Enseco  
Denis Lin, Enseco  
Paul Taylor, Enseco  
Gary Ward, Enseco

SF 5011 MPCA

Haddr G.W.

U.S.E.P.A. - C.L.P.

PAGE : 1

S A M P L E M A N A G E M E N T O F F I C E

INORGANICS  
RESOLUTION OF CONTRACT COMPLIANCE SCREENING (CCS) RESULTS

LABORATORY NAME : ENSECO RMAL

CASE : 09288

RESPONSE RECEIPT DATE : 05/19/88

REGION : 5

SDG\_NO.: MEW615

RECONCILIATION DATE : 05/24/88

RECONCILED BY : JD

DATE MAILED : 05/27/88

ATTACHED ARE COPIES OF CCS SUMMARIES WHICH SHOW THE STATUS OF RELEVANT SAMPLES AFTER INCORPORATION OF  
LABORATORY RESPONSE TO SCREENING. PROBLEM CODES WHICH NO LONGER APPLY ARE MARKED WITH AN (X) CODE .

CRITERION	COMMENTS
B-H	ACTION REMAINS RS.
J	ACTION REMAINS RS.
J2	ACTION REMAINS RS.
R	ACTION REMAINS RS.
I	ACTION REMAINS RS.
	REPORTING OF MSA'S ARE RAS.
R2	ACTION REMAINS RS. SEE ABOVE.

| RECONCILED STATUS OF CCS RESULTS |

PAGE: 2

CONTRACT COMPLIANCE SCREENING SUMMARY FOR INORGANICS

CASE: 09288 CONTRACT: 68-01-7476 SAMPLES: 9 SDG\_NO.: MEW615 DATE SCREENED: 04/29/88

LAB NAME: ENSECO RMAL REGION: 5 SCREENED BY: SW

PROBLEM CODE: E = EXPLAIN, R = RESUBMIT, S = SUBMIT, N = NONCOMPLIANT, X = RECONCILED

PAGE

CONTRACT COMPLIANCE SCREENING SUMMARY FOR INORGANIC S

MPCA

SF 5011

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PRINTED IN U.S.A.

JUN 20. 22

T-10-A. Ground Water  
Water Div.

CASE 09288

SAMPLES · 9

DATE MAILED 03/06/88

LAB NAME ENSECO EMAIL

CONTRACT 68-01-7474

SCREENER S

DATE SCREENED 04/29/88

SDG NUMBER. MEW613

REGION 5

DATE RECEIVED 04/23/88

**PROBLEM CODE E = EXPLAIN, R = RESUBMIT, S = SUBMIT, N = NONCOMPLIANT**

## CONTRACT COMPLIANCE SCREENING FOR INORGANICS

WORKSHEET A - ICP

PAGE: 2

CASE: 09288

CONTRACT: 48-01-7476

SAMPLES: 9

SDQ\_NO.: MEW615

DATE: 4/29/88

LAB NAME: ENSECO /LMAC

REGION 5

SCREENED BY: (R)

(S)

SAMPLE NO.	C/MISAMP	WT.	RECD.	DATE ANAL.	GC ANALYSIS TYPE	TIME	PROBLEMS
MEW615-4	ILW 10.85	02APR88		4/15			(B-H) Sb and Cd values are not reported on standard EPA forms (SOW 787, A-4)
MEW615-1	ILW 11.00	02APR88		12:07	Standard ICV	7:16 - 7:22	(R) 'B-H R
					ICB	7:26	(7) submit traffic reports for all samples. (None were submitted for this case) (SOW 787, B-4)
					CR1	7:31	(S) 'B-H RT
					ICSA	7:34	
					ICSAB	7:39	(FB) Under method, Sb and Cd should have NR as well as Zn. (SOW 787, B-24)
MEW615-9	ILW 11.00	02APR88		12:13	CCVI	11:53	(R) 'B-H FB
					CCBI	11:56	
MEW616-1	ILW 11.00	02APR88		12:17	pBk	11:59	(J) Under method column, enter 'F' for furnace and 'CV' for Hg (SOW 787, B-28)
					CCS(w)	12:03	(e) 'B-H RT
MEW616-8	ILW 11.00	02APR88		12:29			(J2) Serial dilution results do not reflect a 5X dilution ie, $AE = 186.5 \times 5 = 932.5$ 'B-H R
MEW617-1	ILW 11.00	02APR88		12:44			# 187 B (p101) (SOW 787, B-28, E-12)
MEW619-1	ILW 11.00	02APR88		12:48			(R) Report the measured value of each analyte from the raw data to the # of decimal places 'B-H RT
MEW619-1	ILW 11.00	02APR88		12:51			Specified in Exhibit B for the following forms: 2, 4, 7 (SOW 787, B-12, B-21, B-25) (R) 'B-H RT

## CONTRACT COMPLIANCE SCREENING FOR INORGANICS

WORKSHEET A - ICP

PAGE: 3

CASE: 0928B

CONTRACT: 6B-01-7476

SAMPLES: 9

SDG\_ND: MEW615

DATE: 7/29/88

LAB NAME: ENSEC'D /emac

REGION: 5

SCREENED BY: (initials)

SAMPLE NO.	IC/M SAMPLE WT.	DATE RECED.	ANAL.	GC ANALYSIS TYPE	TIME	PROBLEMS
MEW620-11 ILW 1.00	02APR88	13:01	4/15			BHT R
MEW621-1 ILW 1.00	02APR88	13:05	4			BHT R
MEW622-1 ILW 1.00	02APR88	13:12	4			BHT R
MEW623-1 ILW 1.00	02APR88	13:16		SEL.DL: 13:22 CCV3 13:26 CCB3 13:31		BHT R (J.J.)
			(f)	{ CRI 13:35 ICSA 13:39 ICSAB 13:43		

CONTRACT COMPLIANCE SCREENING FOR INORGANICS WORKSHEET B - AA AND CN

PAGE: 4

CASE • 0281

CONTRACT: 68-01-7476

**SAMPLES: Q**

SDG ID : NEW615

DATE: 4/29/88

LAB NAME: ENSECD /emac

RECD ON: 10/10/2018

SCREENED BY: *PJ*

CONTRACT CDMPLIANCE SCREENING FDR INORGANICS WORKSHEET B - AA AND CN

CASE: 09288

**CONTRACT: 6B-01-7476**

SAMPLES: 9

SDG\_ND. : ME4619

**WORKSHEET B - AA AND CNE**

PAGE: 5

LAB NAME: ENSEC'D Lemar

REGION - 5

DATE: 4/29/88

SCREENED BY: *RJ*

SF5211 MPCA

MPCA

**U.S. ENVIRONMENTAL PROTECTION AGENCY  
CONTRACT LABORATORY PROGRAM**

PAGE : 1

# RECEIVED

JUN 24, 88

**ICPCA, Ground Water  
& Solid Waste Div.**

# SAMPLE MANAGEMENT OFFICE

## **RESOLUTION OF CONTRACT COMPLIANCE SCREENING (CCS) RESULT**

**LABORATORY NAME : S3**

CASE NUMBER : 09288

**REGION : 5**

MAIL DATE: 27MAY88

**IDENT : ET831-02**

ATTACHED ARE COPIES OF RECONCILED STATUS OF CCS RESULTS WHICH SHOW THE STATUS OF THE RELEVANT SAMPLES AFTER INCORPORATION OF LABORATORY RESPONSE TO SCREENING. PROBLEM CODES WHICH NO LONGER APPLY ARE MARKED WITH AN "X" CODE.

**COMMENTS : NONE.**

P.O. BOX 818, ALEXANDRIA, VIRGINIA 22313. PHONE : (703) 557-2490/FTS-8-557-2490

## **ATTACHMENTS**

\*\*\*\*\* RECONCILED STATUS OF CCS RESULTS \*\*\*\*\*

PAGE: 2

\*\*\*\*\* CONTRACT COMPLIANCE SCREENING SUMMARY FOR ORGANICS \*\*\*\*\*

**CASE: 09288**      **CONTRACT: 68-01-7261**

**SAMPLES: 12 +**

DATE SCREENED: 04MAY88

**LAB NAME: S3**

**REGION:** 5

**SCREENED BY: SR/AH**

V O A

BNA

PESTICIDES

PROBLEM CODE: E = EXPLAIN, R = RESUBMIT, S = SUBMIT, N = NONCOMPLIANT, NA = FRACTION NOT ANALYZED, X = RECONCILED

**RECEIVED**

**MAY 04 88**

DATE: May 2, 1988

**MPCA, Ground Water  
& Solid Waste Div.**

TO: Minnesota Pollution Control Agency  
Site Assessment Unit  
Program Development Section  
Groundwater and Solid Waste Division  
520 Layfayette Road  
St. Paul, MN 55155

ATTN: Doug Day, Supervisor

Case No.	Site	Contract Lab	SF No.	No. Samples
9288SAS3554E	Hader Ground Water Contamination	RMAL	SF5011	9

Enclosed is one reviewed case.

FROM: U.S. EPA  
Region V  
Central Regional Laboratory  
536 S. Clark, 10th Floor  
Chicago, IL 60605

SENT BY: *Erlinda Day M. Arreola*  
*WESTON/ESAT*

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

DATE: 4.29.88

SUBJECT: Review of Region V CLP Data  
Received for Review on 4-25-88FROM: Curtis Ross, Director (5SCR) *alda leamy*  
Central Regional LaboratoryTO: Data User: MPCA  
Dave Koubeky

We have reviewed the data for the following case(s).

SITE NAME: Hader Ground Water Contamination SMO Case No. 92885AS3554E  
EPA Data Set No. SF 5011 No. of Samples: 9 D.U./Activity Numbers Y905/C 72222CRL No. 88YL06S01-S08, DOSSMO Traffic No. MEW615-623CLP Laboratory: R.MALHrs. Required for Review: 5.5

Following are our findings:

The laboratory's portion of case 9288/SAS3554E consisted of 9 drinking water samples analyzed for total metals and cyanide.

ICP Analysis: Cr matrix spike %R is (84%). Cr data is estimated (UJ) for all samples except sample MEW621, which is estimated (J). Duplicate RPD for Cu (17.4 ug/L) indicate poor precision. Cu is estimated (J).

GFAA Analysis: Matrix spike %R for Cd (80%), Se(205%) and Tl (122.5%) are beyond the +/-15% limit. Se and Tl data are acceptable. Cd data are estimated (UJ) for samples MEW617, 620 and estimated (J) for all the rest. Hg and cyanide data are within control limits.

- (X) Data are acceptable for use.
- ( ) Data are acceptable for use with qualifications referenced above. See Data Qualifier sheets and Calibration Outlier forms for flags and additional comments.
- ( ) Data are preliminary - pending verification by Contractor Laboratory. See Case Summary above.
- ( ) Data are unacceptable.

cc: Carla Dempsey, CLP Quality Assurance Officer, Analytical Operations Branch  
James Petty, Chief Quality Assurance Research, EMSL, Las Vegas*Margaret Ogle**4/28/88*  
*Weston P*

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

ESD/Central Regional Laboratory  
DATA TRACKING FORM FOR CONTRACT SAMPLES

RL Data Set No. SF5011 CERCLIS No. \_\_\_\_\_

SMO Case No. 92885AS 355YE Site Name and Location: Hader Ground Water

Name of Contractor or EPA Laboratory: RMAL Data User: MPCA

No. of Samples: 9 Date Samples or Data Received: 4-25-88

1. Have chain-of-custody records been received? YES  NO
2. Have Traffic Reports or packing lists been received? YES  NO
3. If no, are Traffic Report or packing list numbers written on the chain-of-custody record? YES  NO
4. If no, which Traffic report or packing list numbers are missing?

\_\_\_\_\_

Are basic data forms in? YES  NO

Number of samples claimed: 9 Number of samples received: 9

Checked by: Wanda Yvette Freeman Date: 4-25-88

Received by Contract Project Management Section: IL Date: 4-25-88

Review Started: 4/27/88 Reviewer Signature: Margaret J. Burke

Total time spent on review: 5.5 Date review completed: 4/28/88

Copied (xeroxed) by: Littie + 1 hr verifier by K. Kuse Date: \_\_\_\_\_

Mailed to Data User by: Erinda Lucy M. Aresta Date: 5/2/88 5/2/88

DATA USERS:

Please fill in the blanks below and return this form to: Sylvia Griffin, Data Management Coordinator, Region V, SSCRCL

Data received by: \_\_\_\_\_ Date: \_\_\_\_\_

Q.A. review received by: \_\_\_\_\_ Date: \_\_\_\_\_

Inorganic Data Complete [ ] . Suitable for Intended Purposes [ ]  [ ] if acceptable.  
Organic Data Complete [ ] . Suitable for Intended Purposes [ ] List problems below.  
Dioxin Data Complete [ ] . Suitable for Intended Purposes [ ]  
SAS Data Complete [ ] . Suitable for Intended Purposes [ ]

See Attached "Missing Data Request Form" [ ]

PROBLEMS: Please indicate reasons (if any) why data are not suitable for your uses.  
Other problems.

\_\_\_\_\_

Received by Data Management Coordinator, CRL for File: Date: \_\_\_\_\_

Signature: \_\_\_\_\_

00001

## COVER PAGE - INORGANIC ANALYSES DATA PACKAGE

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615SOW No.: 7/87

## EPA Sample No.

MEW615  
MEW615D  
MEW616  
MEW616A  
MEW616S  
MEW617  
MEW618  
MEW619  
MEW620  
MEW621  
MEW622  
MEW623

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## Lab Sample ID

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U.S. EPA CENTRAL  
REGIONAL LAB

Were ICP interelement corrections applied?

Yes/No YES

Were ICP background corrections applied?

Yes/No YESIf yes-were raw data generated before  
application of background corrections?Yes/No NO

## Comments:

9 LOW WATERS FOR TOTAL METALS AND CYANIDE ANALYSIS  
ALL FURNACE VALUES ARE DETERMINED BY MSA  
RMA QC# 87746 21 DAY TURNAROUND

Release of the data contained in this hardcopy data package and in the  
computer readable data submitted on floppy diskette have been authorized by  
the Laboratory Manager or the Manager's designee, as verified by the  
following signature.

Lab Manager: [Signature]Date: 04/22/88

00002

U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MEW615

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Recieved: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	633	-		P
7440-36-0	Antimony				
7440-38-2	Arsenic	1.7	B	S	F
7440-39-3	Barium	34.4	B		P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium				
7440-70-2	Calcium	167000	-		P
7440-47-3	Chromium	4.0	U	N	P
7440-48-4	Cobalt	16.8	-		P
7440-50-8	Copper	27.4	-	*	P
7439-89-6	Iron	78800	-		P
7439-92-1	Lead	5.5	-	S	F
7439-95-4	Magnesium	47400	-		P
7439-96-5	Manganese	570	-		P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	86.2	-		P
7440-09-7	Potassium	1320	B		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	3570	-		P
7440-28-0	Thallium	2.0	U	SN	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	1340	U		P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: YELLOWClarity Before: CLEAR  
Clarity After: CLEARTexture: \_\_\_\_\_  
Artifacts: \_\_\_\_\_

Comments:

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MEW615

00003

## INORGANIC ANALYSIS DATA SHEET

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	0.3	B	SN	F

Color Before: COLORLESS Clarity Before: CLEAR Texture: \_\_\_\_\_Color After: YELLOW Clarity After: CLEAR Artifacts: \_\_\_\_\_

COMMENTS:

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U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET00004  
EPA SAMPLE NO.Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW616Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Recieved: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	25.0	U		P
7440-36-0	Antimony		U		
7440-38-2	Arsenic	1.0	U	S	F
7440-39-3	Barium	88.2	U		P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium		-		
7440-70-2	Calcium	150000	-		P
7440-47-3	Chromium	4.0	U	N	P
7440-48-4	Cobalt	9.0	B		P
7440-50-8	Copper	10.0	U	*	P
7439-89-6	Iron	5790	-		P
7439-92-1	Lead	3.4	-	S	F
7439-95-4	Magnesium	40200	-		P
7439-96-5	Manganese	181			P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	42.5			P
7440-09-7	Potassium	1300	B		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	2290			P
7440-28-0	Thallium	2.0	U	SN	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	908	U		P
	Cyanide	10.0			AS

Color Before: COLORLESS  
Color After: COLORLESSClarity Before: CLEAR  
Clarity After: CLEARTexture: \_\_\_\_\_  
Artifacts: \_\_\_\_\_

Comments:

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U.S. EPA - CLP

1

EPA SAMPLE NO

MEW616

00005

## INORGANIC ANALYSIS DATA SHEET

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	0.2	B	SN	F

Color Before: COLORLESS Clarity Before: CLEAR Texture: \_\_\_\_\_Color After: COLORLESS Clarity After: CLEAR Artifacts: \_\_\_\_\_

COMMENTS:

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1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO. 00006

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW617Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	25.9	B		P
7440-36-0	Antimony				
7440-38-2	Arsenic	1.0	U	S	F
7440-39-3	Barium	72.6	U		P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium		-		
7440-70-2	Calcium	74300	U		P
7440-47-3	Chromium	4.0	U	N	P
7440-48-4	Cobalt	4.0	U		P
7440-50-8	Copper	10.0	U	*	P
7439-89-6	Iron	1560	-		P
7439-92-1	Lead	3.8	-	S	F
7439-95-4	Magnesium	23200	-		P
7439-96-5	Manganese	68.8	-		P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	10.0	U		P
7440-09-7	Potassium	1030	B		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	3420	U		P
7440-28-0	Thallium	2.0	U	S/N	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	39.1	U		P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: COLORLESSClarity Before: CLEAR  
Clarity After: CLEARTexture:  
Artifacts: \_\_\_\_\_

Comments:

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MEW617

00007

## INORGANIC ANALYSIS DATA SHEET

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID:

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	0.1	U	SN	F

Color Before: COLORLESS Clarity Before: CLEAR Texture: \_\_\_\_\_Color After: COLORLESS Clarity After: CLEAR Artifacts: \_\_\_\_\_

COMMENTS:

1  
INORGANIC ANALYSIS DATA SHEETLab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW618Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	1660	-		P
7440-36-0	Antimony		-		
7440-38-2	Arsenic	3.6	B	S	F
7440-39-3	Barium	91.5			P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium		-		
7440-70-2	Calcium	80000	U	N	P
7440-47-3	Chromium	4.0	U		P
7440-48-4	Cobalt	47.6	U		P
7440-50-8	Copper	10.0	U	*	P
7439-89-6	Iron	83200	-		P
7439-92-1	Lead	9.5	-	S	F
7439-95-4	Magnesium	20700	-		P
7439-96-5	Manganese	655			P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	197	-		P
7440-09-7	Potassium	833	B		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	1800			P
7440-28-0	Thallium	2.0	U	SN	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	1470	U		P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: YELLOWClarity Before: CLEAR  
Clarity After: CLEARTexture:  
Artifacts: \_\_\_\_\_

Comments:

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MEW618

00003

## INORGANIC ANALYSIS DATA SHEET

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID:

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	0.5	B	SN	F

Color Before: COLORLESS Clarity Before: CLEAR Texture: \_\_\_\_\_Color After: YELLOW Clarity After: CLEAR Artifacts: \_\_\_\_\_

COMMENTS:

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00010

EPA SAMPLE NO.

1  
INORGANIC ANALYSIS DATA SHEETLab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW619Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	34.0	B		P
7440-36-0	Antimony		U	S	F
7440-38-2	Arsenic	1.0	B		P
7440-39-3	Barium	28.6	U		P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium		-		
7440-70-2	Calcium	111000			P
7440-47-3	Chromium	4.0	U	N	P
7440-48-4	Cobalt	4.0	U		P
7440-50-8	Copper	10.0	U	*	P
7439-89-6	Iron	9780	U		P
7439-92-1	Lead	2.0	U	S	F
7439-95-4	Magnesium	40600	-		P
7439-96-5	Manganese	621			P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	10.0	U		P
7440-09-7	Potassium	1780	B		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	3230			P
7440-28-0	Thallium	2.0	U	SN	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	32.3	U		P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: YELLOWClarity Before: CLEAR  
Clarity After: CLEARTexture:  
Artifacts: \_\_\_\_\_

Comments:

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MEW619

## INORGANIC ANALYSIS DATA SHEET

00011

Lab Name: ROCKY MOUNTAIN ANALYTICAL

Contract: 68-01-7476

Lab Code: ENSECO

Case No.: 9288

SAS No.: 3554E

SDG No.: MEW615

Matrix (soil/water): WATER

Lab Sample ID:

Level (low/med): LOW

Date Received: 04/02/88

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	0.1	B	SN	F

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: YELLOW

Clarity After: CLEAR

Artifacts:

COMMENTS:

1  
INORGANIC ANALYSIS DATA SHEETLab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW620Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	34.3	B		P
7440-36-0	Antimony				
7440-38-2	Arsenic	1.0	U	S	F
7440-39-3	Barium	28.8	B		P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium				
7440-70-2	Calcium	113000	-		P
7440-47-3	Chromium	4.0	U	N	P
7440-48-4	Cobalt	4.0	U		P
7440-50-8	Copper	10.0	U	*	P
7439-89-6	Iron	9700			P
7439-92-1	Lead	3.4	-	S	F
7439-95-4	Magnesium	41000	-		P
7439-96-5	Manganese	574			P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	10.0	U		P
7440-09-7	Potassium	1780	B		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	2890	-		P
7440-28-0	Thallium	2.0	U	S/N	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	28.5	U		P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: YELLOWClarity Before: CLEAR  
Clarity After: CLEARTexture: \_\_\_\_\_  
Artifacts: \_\_\_\_\_

Comments:

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MEW620

00013

## INORGANIC ANALYSIS DATA SHEET

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID:

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	3.5	B	S	F
7440-43-9	Cadmium	0.1	U	SN	F

Color Before: COLORLESSClarity Before: CLEAR

Texture: \_\_\_\_\_

Color After: YELLOWClarity After: CLEAR

Artifacts: \_\_\_\_\_

COMMENTS:

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1  
INORGANIC ANALYSIS DATA SHEETLab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW621Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Recieved: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	32100	-		P
7440-36-0	Antimony		-		
7440-38-2	Arsenic	8.4	-	S	F
7440-39-3	Barium	55.1	-		P
7440-41-7	Beryllium	4.5	B		P
7440-43-9	Cadmium		-		
7440-70-2	Calcium	92200	-		P
7440-47-3	Chromium	11.0	-	N	P
7440-48-4	Cobalt	113	-		P
7440-50-8	Copper	22.0	-	*	P
7439-89-6	Iron	236000	-		P
7439-92-1	Lead	19.9	-	S	F
7439-95-4	Magnesium	24400	-		P
7439-96-5	Manganese	1670	-		P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	434	-		P
7440-09-7	Potassium	2170	-		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	2810	-		P
7440-28-0	Thallium	2.0	U	SN	F
7440-62-2	Vanadium	25.8	-		P
7440-66-6	Zinc	4120	-		P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: YELLOWClarity Before: CLEAR  
Clarity After: CLEARTexture:  
Artifacts: \_\_\_\_\_

Comments:

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MEW621

## INORGANIC ANALYSIS DATA SHEET

00015

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID:

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	10.0		SN	F

Color Before: COLORLESSClarity Before: CLEAR

Texture:

Color After: YELLOWClarity After: CLEAR

Artifacts:

COMMENTS:

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW622Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Recieved: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	25.0	U		P
7440-36-0	Antimony				
7440-38-2	Arsenic	3.6	B	S	F
7440-39-3	Barium	85.1			P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium				
7440-70-2	Calcium	84000			P
7440-47-3	Chromium	4.0	U	N	P
7440-48-4	Cobalt	4.0	U		P
7440-50-8	Copper	10.0	U	*	P
7439-89-6	Iron	3560			P
7439-92-1	Lead	2.0	U	S	F
7439-95-4	Magnesium	28500			P
7439-96-5	Manganese	69.8			P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	10.0	U		P
7440-09-7	Potassium	5660			P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	6680			P
7440-28-0	Thallium	2.0	U	SN	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	2720	U		P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: COLORLESSClarity Before: CLEAR  
Clarity After: CLEARTexture: \_\_\_\_\_  
Artifacts: \_\_\_\_\_

Comments:

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U.S. EPA - CLP

1

EPA SAMPLE NO

MEW622

## INORGANIC ANALYSIS DATA SHEET

00017

Lab Name: ROCKY MOUNTAIN ANALYTICAL

Contract: 68-01-7476

Lab Code: ENSECO

Case No.: 9288

SAS No.: 3554E

SDG No.: MEW615

Matrix (soil/water): WATER

Lab Sample ID:

Level (low/med): LOW

Date Received: 04/02/88

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	0.1	B	SN	F

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

COMMENTS:

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## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET00018  
EPA SAMPLE NO.Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW623Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID: \_\_\_\_\_

Level (low/med): LOWDate Recieved: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	216	-		P
7440-36-0	Antimony		-		
7440-38-2	Arsenic	5.0	-	S	F
7440-39-3	Barium	90.6			P
7440-41-7	Beryllium	2.0	U		P
7440-43-9	Cadmium		-		
7440-70-2	Calcium	85500			P
7440-47-3	Chromium	4.0	U	N	P
7440-48-4	Cobalt	120			P
7440-50-8	Copper	10.0	U	*	P
7439-89-6	Iron	76800			P
7439-92-1	Lead	2.0	U	S	F
7439-95-4	Magnesium	27000	-		P
7439-96-5	Manganese	517			P
7439-97-6	Mercury	0.2	U		CV
7440-02-0	Nickel	342	-		P
7440-09-7	Potassium	1160	B		P
7482-49-2	Selenium	1.0	U	SN	F
7440-22-4	Silver	4.0	U		P
7440-23-5	Sodium	3600			P
7440-28-0	Thallium	2.0	U	S/N	F
7440-62-2	Vanadium	4.0	U		P
7440-66-6	Zinc	477			P
	Cyanide	10.0	U		AS

Color Before: COLORLESS  
Color After: YELLOWClarity Before: CLEAR  
Clarity After: CLEARTexture: \_\_\_\_\_  
Artifacts: \_\_\_\_\_

Comments:

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MEW623  
00019

## INORGANIC ANALYSIS DATA SHEET

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Matrix (soil/water): WATER

Lab Sample ID:

Level (low/med): LOWDate Received: 04/02/88% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight) : UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-36-0	Antimony	2.0	U	S	F
7440-43-9	Cadmium	0.3	B	SN	F

Color Before: COLORLESSClarity Before: CLEAR

Texture: \_\_\_\_\_

Color After: YELLOWClarity After: CLEAR

Artifacts: \_\_\_\_\_

COMMENTS:  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

# QC EXCEPTION SUMMARY REPORT

CASE # 9288/SAS 3554E SITE Hader G-W Cont.  
 DATA SET # SF5011 LAB RMAI  
 LAB Q.C. # 87746 REVIEWED BY MMT  
 DATE: 4/28/88

MATRIX: water WATER SAMPLE SPK.  
 CONC.: low WATER SAMPLE DUP.  
 MATRIX:            SOIL SAMPLE SPK.  
 CONC.:            SOIL SAMPLE DUP.

	OVERALL CASE QC								MATRIX SPECIFIC QC						SAMPLE SPECIFIC QC		FIELD QC			REGIONAL QC			OTHER / COMMENTS		
	Molding Time	Cal Blanks	Int Calver	Concen Calver	Prep BD AQ	Prep BSL SOL	ICS %R	ICS %	AQ	SOL	Sol Dup RPD	Sol Spk %R	AQ Dup RPD	AQ Spk %R	Ser Diln	AQ	SOL	GFAA Dup	GFAA Spke	Blank	Dup RPD	Spke %R	Blnd Blank	Blnd Spke %R	Spkt Sample RPD
Aluminum																									
Antimony																									
Arsenic																									
Barium																									
Beryllium																									
Cadmium																									
Calcium																									
Chromium																									
Cobalt																									
Copper																									
Iron																									
Lead																									
Magnesium																									
Manganese																									
Mercury																									
Nickel																									
Potassium																									
Selenium																									
Silver																									
Sodium																									
Thallium																									
Tin																									
Vanadium																									
Zinc																									
Cyanide																									

80

84

17.4 >  $\pm 10$  (DL)

205

19.0 (Not flagged on Form I)

122.5

## U.S. EPA - CLP

2A

## INITIAL AND CONTINUING CALIBRATION VERIFICATION

00020

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Initial Calibration Source: EPA-LVContinuing Calibration Source: FISHER

Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration				M
	True	Found	%R(1)	True	Found	%R(1)	Found	
Aluminum	1980.0	2060	104.0	1000.0	977	97.7	998	99.8
Antimony								P
Arsenic								
Barium	1980.0	2020	102.0	1000.0	985	98.5	1010	101.0
Beryllium	481.0	493	102.5	1000.0	988	98.8	1000	100.0
Cadmium								
Calcium	49800.0	53900	108.2	100000.	103000	103.0	105000	105.0
Chromium	506.0	462	91.3	1000.0	913	91.3	936	93.6
Cobalt	474.0	516	108.9	1000.0	1000	100.0	1020	102.0
Copper	542.0	570	105.2	1000.0	1030	103.0	1050	105.0
Iron	1990.0	2060	103.5	1000.0	1010	101.0	1030	103.0
Lead								
Magnesium	25000.0	25900	103.6	100000.	102000	102.0	106000	106.0
Manganese	513.0	540	105.3	1000.0	1040	104.0	1070	107.0
Mercury								
Nickel	496.0	503	101.4	1000.0	1040	104.0	1080	108.0
Potassium	50200.0	51300	102.2	100000.	97500	97.5	99800	99.8
Selenium								
Silver	509.0	480	94.3	1000.0	924	92.4	952	95.2
Sodium	50700.0	52700	103.9	100000.	100000	100.0	104000	104.0
Thallium								
Vanadium	511.0	513	100.4	1000.0	995	99.5	1020	102.0
Zinc	3100.0	3180	102.6	1000.0	1020	102.0	1050	105.0
Cyanide								

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

2A

## INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Initial Calibration Source: EPA-LVContinuing Calibration Source: FISHER

Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration				M
	True	Found	%R(1)	True	Found	%R(1)	Found	
Aluminum				1000.0	994	99.4		P
Antimony								
Arsenic								
Barium				1000.0	1000	100.0		P
Beryllium				1000.0	996	99.6		P
Cadmium								
Calcium				100000.	104000	104.0		P
Chromium				1000.0	925	92.5		P
Cobalt				1000.0	1020	102.0		P
Copper				1000.0	1040	104.0		P
Iron				1000.0	1020	102.0		P
Lead								
Magnesium				100000.	104000	104.0		P
Manganese				1000.0	1060	106.0		P
Mercury								
Nickel				1000.0	1060	106.0		P
Potassium				100000.	98400	98.4		P
Selenium								
Silver				1000.0	943	94.3		P
Sodium				100000.	102000	102.0		P
Thallium								
Vanadium				1000.0	1010	101.0		P
Zinc				1000.0	1040	104.0		P
Cyanide								

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

2A  
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476

Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615

Initial Calibration Source: EPA-LV

Continuing Calibration Source: FISHER

Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration			M
	True	Found	%R(1)	True	Found	%R(1)	
Aluminum							
Antimony							
Arsenic	47.0	45.1	96.0	20.0	19.1	95.5	20.0
Barium							
Beryllium							
Cadmium							
Calcium							
Chromium							
Cobalt							
Copper							
Iron							
Lead	97.9	105	107.3	20.0	20.4	102.0	20.2
Magnesium							
Manganese							
Mercury	5.2	5.6	107.7	1.0	1.0	100.0	1.0
Nickel							
Potassium							
Selenium	104.0	104	100.0	20.0	22.0	110.0	19.6
Silver							
Sodium							
Thallium	97.3	106	108.9	20.0	18.9	94.5	20.8
Vanadium							
Zinc							
Cyanide	105.0	101	96.2	200.0	204	102.0	208
							AS

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2A  
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476  
 Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615  
 Initial Calibration Source: EPA-LV  
 Continuing Calibration Source: FISHER

Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration				M
	True	Found	%R(1)	True	Found	%R(1)	Found	
Aluminum								
Antimony								
Arsenic				20.0	20.3	101.5	20.4	102.0
Barium								
Beryllium								
Cadmium								
Calcium								
Chromium								
Cobalt								
Copper								
Iron								
Lead				20.0	19.8	99.0	19.6	98.0
Magnesium								
Manganese								
Mercury				1.0	1.0	100.0		
Nickel								
Potassium								
Selenium				20.0	20.8	104.0	21.0	105.0
Silver								
Sodium								
Thallium				20.0	19.3	96.5	19.8	99.0
Vanadium								
Zinc								
Cyanide				200.0	208	104.0		

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

2A  
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Initial Calibration Source: EPA-LVContinuing Calibration Source: FISHER

Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration			M
	True	Found	%R(1)	True	Found	%R(1)	
Aluminum							
Antimony							
Arsenic				20.0	21.9	109.5	
Barium							
Beryllium							
Cadmium							
Calcium							
Chromium							
Cobalt							
Copper							
Iron							
Lead							
Magnesium							
Manganese							
Mercury							
Nickel							
Potassium				20.0	21.0	105.0	
Selenium							
Silver							
Sodium							
Thallium							
Vanadium							
Zinc							
Cyanide							

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

<sup>2A</sup>  
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Initial Calibration Source: EPA-LVContinuing Calibration Source: FISHER

Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration				M
	True	Found	%R(1)	True	Found	%R(1)	Found	
Aluminum								
Antimony								
Arsenic								
Barium								
Beryllium								
Cadmium								
Calcium								
Chromium								
Cobalt								
Copper								
Iron								
Lead	97.9	105	107.3	20.0	20.0	100.0	19.7	98.5
Magnesium								
Manganese								
Mercury								
Nickel								
Potassium								
Selenium	104.0	107	102.9	20.0	20.9	104.5	21.0	105.0
Silver								
Sodium								
Thallium	97.3	87.4	89.8	20.0	20.3	101.5	18.0	90.0
Vanadium								
Zinc								
Cyanide								

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP  
2A  
INITIAL AND CONTINUING CALIBRATION VERIFICATION

00026

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476  
 Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615

Initial Calibration Source: EPA-LV

Continuing Calibration Source: FISHER

Concentration Units:ug/L

Analyte	Initial Calibration			Continuing Calibration					M
	True	Found	%R(1)	True	Found	%R(1)	Found	%R(1)	
Antimony	1010.0	940	93.1	20.0	20.6	103.0	20.0	100.0	F
Cadmium	96.1	90.0	93.6	2.0	2.0	100.0	2.1	105.0	F

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): UG/L

Analyte	Initial Calib. Blank (ug/L)	Continuing Calibration Blank						Prepa- ration Blank	C	M
		1	C	2	C	3	C			
Antimony	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U
Cadmium	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP  
2A  
INITIAL AND CONTINUING CALIBRATION VERIFICATION

00027

Lab Name: ROCKY MOUNTAIN ANALYTICAL

Contract: 68-01-7476

Lab Code: ENSECO Case No.: 9288

SAS No.: 3554E SDG No.: MEW615

Initial Calibration Source: EPA-LV

Continuing Calibration Source: FISHER

Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration					M
	True	Found	%R(1)	True	Found	%R(1)	Found	%R(1)	
Antimony				20.0	17.9	89.5	18.5	92.5	F
Cadmium				2.0	2.2	110.0	1.8	90.0	F

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): UG/L

Analyte	Initial Calib. Blank (ug/L)	Continuing Calibration Blank						Prepa- ration Blank	C	M
		1	C	2	C	3	C			
Antimony		2.0	U							F
Cadmium		0.1	U							F

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP  
2A  
INITIAL AND CONTINUING CALIBRATION VERIFICATION

00028

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476  
 Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615

Initial Calibration Source: EPA-LV  
 Continuing Calibration Source: FISHER

Concentration Units:ug/L

Analyte	Initial Calibration			Continuing Calibration				M
	True	Found	%R(1)	True	Found	%R(1)	Found	
Antimony								
Cadmium	96.1	95.0	98.9	2.0	1.9	95.0	1.8	90.0 F

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): UG/L

Analyte	Initial Calib. Blank (ug/L)	Continuing Calibration Blank						Prepa- ration Blank C	M
		1	C	2	C	3	C		
Antimony									
Cadmium	-0.1	B	0.1	U	0.1	U			F

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2B

00029

Lab Name: ROCKY MOUNTAIN ANALYTICAL

**Contract:** 68-01-7476

**Lab Code:** ENSECO      **Case No.:** 9288

SAS No.: 3554E SDG No.: MEW615

AA CRDL Standard Source: BAKER

## ICP CRDL Standard Source: BAKER

**Concentration Units: ug/L**

U.S. EPA - CLP

60030

2B  
CRDL STANDARD FOR AA AND ICP

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECO Case No.: 9288SAS No.: 3554E SDG No.: MEW615AA CRDL Standard Source: BAKERICP CRDL Standard Source: BAKER

Concentration Units: ug/L

Analyte	CRDL Standard for AA			CRDL Standard for ICP		
	True	Found	%R	Initial	Found	Final
Aluminum						
Antimony						
Arsenic						
Barium						
Beryllium						
Cadmium						
Calcium						
Chromium						
Cobalt						
Copper						
Iron						
Lead	2.0	2.4	120.0			
Magnesium						
Manganese						
Mercury						
Nickel						
Potassium						
Selenium	2.0	1.4	70.0			
Silver						
Sodium						
Thallium	1.0	2.4	0.0			
Vanadium						
Zinc						

00031

<sup>2B</sup>  
CRDL STANDARD FOR AA AND ICP

Lab Name: ROCKY MOUNTAIN ANALYTICAL

Contract: 68-01-7476

Lab Code: ENSECO Case No.: 9288

SAS No.: 3554E SDG No.: MEW615

AA CRDL Standard Source: BAKER

ICP CRDL Standard Source: BAKER

Concentration Units: ug/L

Analyte	CRDL Standard for AA			CRDL Standard for ICP			
	True	Found	%R	True	Found	%R	Found
Antimony	5.0	4.2	84.0				

3  
BLANKS

00032

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Preparation Blank Matrix (soil/water): WATERPreparation Blank Concentration Units (ug/L or mg/kg): UG/L

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum	25.0	U	25.0	U	25.0	U	25.0	U	25.0	U	P
Antimony											
Arsenic											
Barium	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U	P
Beryllium	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U	P
Cadmium											
Calcium	81.0	U	81.0	U	81.0	U	81.0	U	81.0	U	P
Chromium	4.0	U	4.0	U	4.0	U	4.0	U	4.0	U	P
Cobalt	4.0	U	4.0	U	4.0	U	4.0	U	4.0	U	P
Copper	10.0	U	10.0	U	10.0	U	10.0	U	10.0	U	P
Iron	14.0	U	14.0	U	-15.0	B	-15.0	B	14.0	U	P
Lead											
Magnesium	81.0	U	81.0	U	81.0	U	81.0	U	81.0	U	P
Manganese	6.0	U	6.0	U	6.0	U	6.0	U	6.0	U	P
Mercury											
Nickel	10.0	U	10.0	U	10.0	U	10.0	U	10.0	U	P
Potassium	95.0	U	95.0	U	95.0	U	95.0	U	95.0	U	P
Selenium											
Silver	4.0	U	4.0	U	4.0	U	4.0	U	4.0	U	P
Sodium	1360	U	1360	U	1360	U	1360	U	1360	U	P
Thallium											
Vanadium	4.0	U	4.0	U	4.0	U	4.0	U	4.0	U	P
Zinc	3.0	U	3.0	U	3.0	U	3.0	U	3.9	B	P
Cyanide											

U.S. EPA - CLP

3  
BLANKS

00033

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Preparation Blank Matrix (soil/water): WATERPreparation Blank Concentration Units (ug/L or mg/kg): UG/L

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum	-	-	-	-	-	-	-	-	-	-	-
Antimony	-	-	-	-	-	-	-	-	-	-	-
Arsenic	1.0	U	1.0	U	1.0	U	1.0	U	1.0	U	F
Barium	-	-	-	-	-	-	-	-	-	-	-
Beryllium	-	-	-	-	-	-	-	-	-	-	-
Cadmium	-	-	-	-	-	-	-	-	-	-	-
Calcium	-	-	-	-	-	-	-	-	-	-	-
Chromium	-	-	-	-	-	-	-	-	-	-	-
Cobalt	-	-	-	-	-	-	-	-	-	-	-
Copper	-	-	-	-	-	-	-	-	-	-	-
Iron	-	-	-	-	-	-	-	-	-	-	-
Lead	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U	F
Magnesium	-	-	-	-	-	-	-	-	-	-	-
Manganese	-	-	-	-	-	-	-	-	-	-	-
Mercury	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	CV
Nickel	-	-	-	-	-	-	-	-	-	-	-
Potassium	-	-	-	-	-	-	-	-	-	-	-
Selenium	1.0	U	1.0	U	1.0	U	1.0	U	1.0	U	F
Silver	-	-	-	-	-	-	-	-	-	-	-
Sodium	-	-	-	-	-	-	-	-	-	-	-
Thallium	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U	F
Vanadium	-	-	-	-	-	-	-	-	-	-	-
Zinc	-	-	-	-	-	-	-	-	-	-	-
Cyanide	10.0	U	10.0	U	10.0	U	10.0	U	10.0	U	AS

3  
BLANKSLab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Preparation Blank Matrix (soil/water): WATERPreparation Blank Concentration Units (ug/L or mg/kg): UG/L

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum											
Antimony											
Arsenic			1.5	B	2.2	B					
Barium											
Beryllium											
Cadmium											
Calcium											
Chromium											
Cobalt											
Copper											
Iron											
Lead			2.0	U							
Magnesium											
Manganese											
Mercury											
Nickel											
Potassium											
Selenium			1.0	U	1.0	U					
Silver											
Sodium											
Thallium			2.0	U							
Vanadium											
Zinc											
Cyanide											

3  
BLANKS

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Preparation Blank Matrix (soil/water): WATERPreparation Blank Concentration Units (ug/L or mg/kg): UG/L

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum											
Antimony											
Arsenic											
Barium											
Beryllium											
Cadmium											
Calcium											
Chromium											
Cobalt											
Copper											
Iron											
Lead	2.0	U	2.0	U	2.0	U					F
Magnesium											
Manganese											
Mercury											
Nickel											
Potassium											
Selenium	1.0	U	1.0	U	1.0	U					F
Silver											
Sodium											
Thallium	2.0	U	2.0	U	2.0	U					F
Vanadium											
Zinc											
Cyanide											

00036

4  
ICP INTERFERENCE CHECK SAMPLELab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9188SAS No.: 3554E SDG No.: new 615ICP ID Number: JA9000ICS Source: EPA-LV-0387

Concentration Units: ug/L

Analyte	True		Initial Found			Final Found		
	Sol.	Sol.	Sol.	Sol.	%R	Sol.	Sol.	%R
	A	AB	A	AB		A	AB	
Aluminum	502000	503000	509000	515000	101.4	531000	514000	101.2
Antimony								
Arsenic								
Barium		483	7	473	97.9	8	585	100.4
Beryllium		474	-0	488	103.0	-0	484	102.1
Cadmium								
Calcium	506000	516000	523000	527000	102.1	537000	526000	101.9
Chromium		513	43	469	91.4	40	469	91.4
Cobalt		478	27	497	104.0	7	516	107.9
Copper		534	18	574	107.5	6	572	107.1
Iron	196000	203000	190000	194000	95.6	216000	194000	95.6
Lead								
Magnesium	497000	509000	517000	524000	102.9	564000	536000	105.3
Manganese		531	93	576	108.5	65	597	112.4
Mercury								
Nickel		916	44	954	104.1	17	975	106.4
Potassium			84	64.5		29	-119	
Selenium								
Silver		993	-0	930	93.7	-724	910	91.6
Sodium			-753	-1160			-625	
Thallium								
Vanadium		415	6	485	102.1	7	487	102.5
Zinc		973	5	984	101.1	-0	1020	104.8

5A  
SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW616SLab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATERLevel (low/med): LOWConcentration Units (ug/L or mg/kg dry weight): UG/L

Analyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Aluminum	75-125	1870	-	25.0	U	2000	93.5	-	P
Antimony									
Arsenic	75-125	19.8	-	1.0	U	20.0	99.0	-	F
Barium	75-125	1930	-	88.2	U	2000	92.1	-	P
Beryllium	75-125	46.5	-	2.0	U	50.0	93.0	-	P
Cadmium									
Calcium	75-125	198000	-	150000	U	50000	96.0	-	P
Chromium	75-125	168	-	4.0	U	200	84.0	N	P
Cobalt	75-125	492	-	9.0	B	500	96.6	-	P
Copper	75-125	253	-	10.0	U	250	101.2	-	P
Iron									
Iron		6320	-	5790	-	1000	53.0	-	P
Lead	75-125	20.4	-	3.4	-	20.0	85.0	-	F
Magnesium	75-125	64800	-	40200	-	25000	98.4	-	P
Manganese	75-125	369	-	181	-	200	94.0	-	P
Mercury	75-125	1.0	-	0.2	U	1.0	100.0	-	CV
Nickel	75-125	426	-	42.5	-	400	95.9	-	P
Potassium	75-125	20200	-	1300	B	20000	94.5	-	P
Selenium	75-125	20.5	-	1.0	U	10.0	205.0	N	F
Silver	75-125	42.8	-	4.0	U	50.0	85.6	-	P
Sodium	75-125	56200	-	2290	-	50000	107.8	-	P
Thallium	75-125	24.5	-	2.0	U	20.0	122.5	N	F
Vanadium	75-125	453	-	4.0	U	500	90.6	-	P
Zinc									
Zinc		1050	-	908	-	200	71.0	-	P
Cyanide	75-125	96.7	-	10.0	U	100	96.7	-	AS

## Comments:

CONTROL LIMIT IS 85 TO 115% FOR ALL METALS EXCEPT MERCURY AND CYANIDE.  
MERCURY AND CYANIDE CONTROL LIMIT IS 80 TO 120% RECOVERY.

U.S. EPA - CLP  
5A  
SPIKE SAMPLE RECOVERY

00038  
EPA SAMPLE NO.

MEW616S

Lab Name: ROCKY MOUNTAIN ANALYTICAL

Contract: 68-01-7476

Lab Code: ENSECO

Case No.: 9288

SAS No.: 3554E

SDG No.: MEW615

Matrix (soil/water): WATER

Level (low/medium): LOW

Concentration Units (ug/L or mg/kg dry weight): UG/L

Analyte	Control Limit %R	Spike Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Antimony	85-115	18.9		2.0	U	20.0	94.5		F
Cadmium	85-115	1.6		0.2	U	2.0	80.0	N	F

EPA SAMPLE NO.

DUPLICATES

MEW615D

% Solids for Sample: 0.0

% Solids for Duplicate: 0.0

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Antimony		2.0	U	2.0	U			F
Cadmium	5.0	0.3	B	0.4	B	28.6		F

Comments:

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00039

EPA SAMPLE NO.

5B  
POST DIGEST SPIKE SAMPLE RECOVERYLab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476mew616ALab Code: ENSECO Case No.: 9288 SAS No.: 35546 SDG No.: mew615Matrix (soil/water): WATERLevel (low/med): LOW

Concentration Units: ug/L

Analyte	Control Limit %R	Spiked Sample Result (SSR) C	Sample Result (SR) C	Spike Added (SA)	%R	Q	M
Aluminum							NR
Antimony							NR
Arsenic							NR
Barium							NR
Beryllium							NR
Cadmium							NR
Calcium							NR
Chromium		19.3	4.0	u 10.0	96.5		NR
Cobalt							NR
Copper							NR
Iron							NR
Lead							NR
Magnesium							NR
Manganese							NR
Mercury							NR
Nickel							NR
Potassium							NR
Selenium							NR
Silver							NR
Sodium							NR
Thallium							NR
Vanadium							NR
Zinc							P
Cyanide							NR

Comments:

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6  
DUPLICATES

EPA SAMPLE NO.

MEW615D

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476Lab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATERLevel (low/med): LOW% Solids for Sample: 0.0% Solids for Duplicate: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Aluminum		633	-	631	-	0.3	-	P
Antimony								
Arsenic		1.7	B	1.1	B	42.9	-	F
Barium		34.4	B	33.5	B	2.7	-	P
Beryllium		2.0	U	2.0	U	-	-	P
Cadmium								
Calcium		167000	U	168000	U	0.6	-	P
Chromium		4.0	U	4.0	U	-	-	P
Cobalt	10.0	16.8	-	15.1	-	10.7	-	P
Copper	10.0	27.4	-	10.0	U	200	*	p
Iron		78800		78800		0.0	-	P
Lead	2.0	5.5	-	3.7	-	39.1	-	F
Magnesium		47400	-	47500	-	0.2	-	P
Manganese		570	-	571	-	0.2	-	P
Mercury		0.2	U	0.2	U	-	-	CV
Nickel	20.0	86.2	U	86.9	U	0.8	-	P
Potassium		1320	B	1330	B	0.8	-	P
Selenium		1.0	U	1.0	U	-	-	F
Silver		4.0	U	4.0	U	-	-	P
Sodium		3570	U	2950	U	19.0	-	P
Thallium		2.0	U	2.0	U	-	-	F
Vanadium		4.0	U	4.0	U	-	-	P
Zinc		1340	U	1340	U	0.0	-	P
Cyanide		10.0	U	10.0	U	-	-	AS

7  
LABORATORY CONTROL SAMPLELab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Solid LCS Source: EMSL - LVAqueous LCS Source: EMSL - LV

Analyte	Aqueous (ug/L)			Solid (mg/kg)			%R
	True	Found	%R(1)	True	Found	C	
Aluminum	1980	1970	99.5			-	
Antimony						-	
Arsenic						-	
Barium	1980	1960	99.0			-	
Beryllium	481	506	105			-	
Cadmium						-	
Calcium	49800	51400	103			-	
Chromium	506	442	87.4			-	
Cobalt	474	504	106			-	
Copper	542	557	103			-	
Iron	1990	1970	99.0			-	
Lead						-	
Magnesium	25000	25100	100			-	
Manganese	513	534	104			-	
Mercury						-	
Nickel	496	490	98.8			-	
Potassium	50200	47100	93.8			-	
Selenium						-	
Silver	509	453	89.0			-	
Sodium	50700	49600	97.8			-	
Thallium						-	
Vanadium	511	489	95.7			-	
Zinc	3100	3060	98.7			-	
Cyanide						-	

7  
LABORATORY CONTROL SAMPLELab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615Solid LCS Source: EMSL - LVAqueous LCS Source: EMSL - LV

Analyte	Aqueous (ug/L)			Solid (mg/kg)				%R
	True	Found	%R(1)	True	Found	C	Limits	
Aluminum								
Antimony								
Arsenic	47.0	42.9	91.3					
Barium								
Beryllium								
Cadmium								
Calcium								
Chromium								
Cobalt								
Copper								
Iron								
Lead	97.9	86.8	88.7					
Magnesium								
Manganese								
Mercury								
Nickel								
Potassium								
Selenium	104	104	100					
Silver								
Sodium								
Thallium	97.3	84.6	86.9					
Vanadium								
Zinc								
Cyanide								

7  
LABORATORY CONTROL SAMPLE

00043

Lab Name: ROCKY MOUNTAIN ANALYTICALContract : 68-01-7476Lab Code: ENSECO Case No.: 9288SAS No.: 3554E SDG No.: MEW615Solid LCS Source: EMSL-LVAqueous LCS Source: EMSL-LV

Analyte	Aqueous (ug/L)			Solid (mg/kg)				
	True	Found	%R	True	Found	C	Limits	%R
Antimony	1010.0	885	87.6					
Cadmium	96.1	90.0	93.6					

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STANDARD ADDITION RESULTS

00044

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615

Concentration Units: ug/L

EPA Sample No.	An	Dil	0 ADD ABS	1 ADD CON ABS	2 ADD CON ABS	3 ADD CON ABS	Final Conc.	r	Q
PB	AS	1	0.000	10	10.000	20	21.700	14 0.0	0.9995
LCS	AS	1	45.800	10	57.000	20	67.600	42.9	0.9999
CCB1	AS	1	0.000	10	10.100	20	19.600	14 0.0	0.9992
CCB2	AS	1	0.000	10	10.500	20	20.000	14 0.0	0.9991
CCB3	AS	1	0.000	10	10.300	20	20.100	30 34.800	0.9952
CCB4	AS	1	2.200	10	12.200	20	22.700	30 35.400	1.5
CCV4	AS	1	23.000	10	34.600	20	43.900	30 57.200	0.9984
MEW621	AS	1	9.200	10	19.200	20	30.500	30 41.000	8.4
MEW616S	AS	1	21.900	10	42.800	20	57.100	30 67.500	16.3
MEW616S	AS	1	22.400	10	36.700	20	46.300	30 58.400	19.8
CCV1	AS	1	19.900	10	28.400	20	40.100	30 49.800	19.1
ICV	AS	1	46.600	10	56.900	20	67.200	30 77.600	45.1
CCV3	SB	1	19.500	10	28.900	20	42.500	30 50.800	17.9
MEW620	AS	1	0.000	10	11.700	20	24.600	30 37.800	14 0.0
CCV3	AS	1	20.000	10	29.900	20	42.000	30 49.300	20.3
CCB3	SB	1	0.000	10	9.800	20	19.300	30 30.100	14 0.0
MEW617	AS	1	0.000	10	11.900	20	23.800	30 35.300	14 0.0
CCV2	AS	1	20.200	10	30.200	20	41.500	30 50.300	20.0
MEW616	AS	1	1.300	10	12.700	20	25.800	30 38.800	14 0.0
CRA	AS	1	4.800	10	15.000	20	25.900	30 35.400	4.7
CCV5	AS	1	22.100	10	30.800	20	42.200	30 51.400	21.9
CCB5	AS	1	2.100	10	12.400	20	23.000	30 32.300	2.2
ICB	SB	1	0.000	10	7.800	20	16.000	30 25.900	14 0.0
LCS	SB	50	19.800	10	30.400	20	42.800	30 52.800	885.0

## STANDARD ADDITION RESULTS

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615

Concentration Units: ug/L

EPA Sample No.	An	Dil	0 ADD ABS	1 ADD CON ABS	2 ADD CON ABS	3 ADD CON ABS	Final Conc.	r	Q
CCB4	SB	1	0.000	10	10.200	20	20.300	30	30.900
ICB	AS	1	0.000	10	9.300	20	20.200	30	31.200
CCV4	SB	1	20.300	10	32.400	20	43.200	30	53.900
CCB2	SB	1	0.000	10	8.700	20	17.600	30	28.200
CCV1	SB	1	19.700	10	29.600	20	39.200	30	48.600
ICV	SB	50	18.300	10	28.400	20	38.100	30	47.700
PB	SB	1	1.800	10	10.200	20	19.100	30	29.800
CCB1	SB	1	0.000	10	8.200	20	18.800	30	27.700
CCV2	SB	1	20.400	10	30.900	20	39.700	30	51.400
CRA	SB	1	4.900	10	13.000	20	23.400	30	34.300
MEW622	AS	1	5.000	10	17.000	20	30.700	30	43.600
MEW623	SB	1	0.000	10	8.400	20	19.200	30	27.800
MEW622	SB	1	0.000	10	10.800	20	24.700	30	39.700
MEW619	SB	1	0.000	10	11.000	20	21.000	30	33.600
MEW618	SB	1	0.000	10	7.500	20	17.000	30	25.500
MEW623	AS	1	5.600	10	16.900	20	28.700	30	39.300
MEW619	AS	1	0.000	10	12.300	20	25.900	30	39.700
MEW615	SB	1	0.000	10	4.800	20	11.000	30	16.500
MEW615	AS	1	2.200	10	11.600	20	23.400	30	33.500
MEW621	SB	1	2.500	10	11.100	20	24.800	30	38.400
MEW616	SB	1	0.000	10	7.800	20	17.500	30	27.600
MEW620	SB	1	3.700	10	15.100	20	24.900	30	36.400
MEW618	AS	1	4.600	10	15.100	20	26.200	30	38.700
MEW617	SB	1	0.000	10	10.500	20	24.700	30	37.800

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## **STANDARD ADDITION RESULTS**

00046

Lab Name: ROCKY MOUNTAIN ANALYTICAL

**Contract:** 68-01-7476

**Lab Code: ENSECO**

**Case No.:** 9288

**SAS No.: 3554E**

**SDG NO.: MEW615**

**Concentration Units: ug/L**

8  
STANDARD ADDITION RESULTS

00047

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554E SDG No.: MEW615

Concentration Units: ug/L

EPA Sample No.	An	Dil	0 ADD ABS	1 ADD CON	1 ADD ABS	2 ADD CON	2 ADD ABS	3 ADD CON	3 ADD ABS	Final Conc.	r	Q
PB	CD	1	0.000	1	0.400	1	1.000	2	1.500	0.140.0	0.9973	-
LCS	CD	50	1.900	1	2.500	1	3.000	2	3.500	90.0	0.9989	-
CCV1	CD	1	1.800	1	2.300	1	2.800	2	3.200	1.9	0.9986	-
CCV2	CD	1	1.800	1	2.300	1	2.800	2	3.300	1.8	1.0000	-
ICV1	CD	50	1.900	1	2.300	1	2.800	2	3.400	90.0	0.9960	-
MEW616S	CD	1	1.700	1	2.100	1	2.600	2	3.200	1.6	0.9960	-
ICB	PB	1	0.000	10	10.200	20	18.700	30	29.400	240.0	0.9991	-
CCB2	PB	1	0.000	10	9.400	20	20.900	30	31.200	240.0	0.9993	-
CCV2	PB	1	19.900	10	28.700	20	38.800	30	49.400	19.7	0.9991	-
ICV	CD	50	1.900	1	2.300	1	2.900	2	3.300	95.0	0.9965	-
CCV1	PB	1	19.400	10	30.100	20	39.100	30	49.200	20.0	0.9995	-
CCV3	PB	1	19.200	10	28.400	20	39.700	30	47.300	20.2	0.9975	-
ICB	CD	1	-0.100	1	0.400	1	0.900	2	1.400	-0.1	1.0000	-
CCV1	CD	1	2.000	1	2.600	1	3.100	2	3.500	2.0	0.9960	-
CCB1	CD	1	0.000	1	0.500	1	1.000	2	1.500	0.140.0	1.0000	-
ICV	PB	1	47.300	10	55.100	20	63.700	30	74.100	52.5	0.9978	-
CRA	PB	1	2.700	5	7.200	10	12.900	15	18.800	2.1	0.9982	-
ICV1	PB	2	49.300	10	57.800	20	66.500	30	77.400	105.0	0.9981	-
ICB1	CD	1	0.000	1	0.500	1	1.000	2	1.500	0.140.0	1.0000	-
CCV4	CD	1	1.900	1	2.500	1	3.100	2	3.500	1.8	0.9959	-
CCB2	CD	1	-0.100	1	0.400	1	0.800	2	1.300	0.140.0	0.9991	-
CCB3	PB	1	0.000	10	9.300	20	20.300	30	28.200	240.0	0.9982	-
CCB1	CD	1	-0.100	1	0.400	1	0.900	2	1.300	0.140.0	0.9986	-
CCB3	CD	1	0.000	1	0.500	1	1.000	2	1.600	0.140.0	0.9989	-

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00048

## STANDARD ADDITION RESULTS

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615

Concentration Units: ug/L

EPA Sample No.	An	Dil	0 ADD ABS	1 ADD CON ABS	2 ADD CON ABS	3 ADD CON ABS	Final Conc.	r	Q
CCV2	CD	1	2.100	1 2.600	1 3.100	2 3.600	2.1	1.0000	-
CCB4	CD	1	0.000	1 0.500	1 1.100	2 1.600	0.1u0.0	0.9993	-
CCB5	PB	1	0.000	10 9.400	20 21.300	30 29.000	0.1u0.0	0.9970	-
CCB1	PB	1	0.000	10 10.200	20 20.400	30 30.700	2u 0.0	1.0000	-
CCB1	PB	1	0.000	10 9.500	20 20.800	30 29.700	2u 0.0	0.9990	-
CCV5	PB	1	19.900	10 29.600	20 40.600	30 49.900	19.6	0.9995	-
PB	PB	1	1.200	10 11.200	20 21.300	30 31.500	1.1	1.0000	-
CCV1	PB	1	20.500	10 30.500	20 40.300	30 50.600	20.4	1.0000	-
CRA	SB	1	4.900	10 13.000	20 23.400	30 34.300	4.2	0.9978	-
CCV4	PB	1	19.900	10 29.400	20 39.700	30 49.700	19.8	0.9999	-
ICB1	PB	1	0.000	10 9.400	20 19.500	30 28.300	2u 0.0	0.9997	-
CRA	PB	1	2.900	5 7.600	10 13.700	15 19.000	2.4	0.9989	-
CCV3	CD	1	2.000	1 2.500	1 2.900	2 3.400	2.2	0.9991	-
CCB4	PB	1	0.000	10 9.600	20 21.200	30 29.000	2u 0.0	0.9974	-
CCB2	CD	1	0.000	1 0.500	1 1.000	2 1.600	0.1u0.0	0.9989	-
LCS1	PB	2	43.000	10 54.200	20 64.300	30 73.100	86.8	0.9986	-
MEW620	CD	1	0.100	1 0.500	1 0.900	2 1.400	0.1u0.0	0.9984	-
MEW622	CD	1	0.100	1 0.500	1 1.000	2 1.300	0.1	0.9959	-
MEW616	CD	1	0.200	1 0.700	1 1.100	2 1.600	0.2	0.9991	-
MEW622	PB	1	0.000	10 9.500	20 18.400	30 26.200	2u 0.0	0.9990	-
MEW621	PB	1	15.200	10 22.300	20 30.000	30 37.800	19.9	0.9998	-
MEW616	PB	1	2.600	10 10.200	20 17.600	30 25.500	3.4	0.9999	-
MEW615	PB	1	4.300	10 11.600	20 20.000	30 27.000	5.5	0.9994	-
MEW617	CD	1	0.100	1 0.500	1 1.000	2 1.500	0.1u0.0	0.9986	-

## 8 STANDARD ADDITION RESULTS

Lab Name: ROCKY MOUNTAIN ANALYTICAL

**Contract:** 68-01-7476

**Lab Code:** ENSECO

**Case No.:** 9288

SAS No.: 3554E

SDG No.: MEW615

**Concentration Units: ug/L**

8  
STANDARD ADDITION RESULTS

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554E SDG No.: MEW615

Concentration Units: ug/L

EPA Sample No.	An	Dil	0 ADD ABS	1 ADD CON ABS	2 ADD CON ABS	3 ADD CON ABS	Final Conc.	r	Q
PB	SE	1	1.100	10 12.700	20 24.900	30 36.300	14 0.0	0.9999	-
CRA	SE	1	1.400	10 10.600	20 23.000	30 32.700	14 0.0	0.9984	-
CCB1	SE	1	0.000	10 10.100	20 19.600	30 31.200	14 0.0	0.9991	-
CCB5	SE	1	0.000	10 12.100	20 23.300	30 33.800	14 0.0	0.9995	-
CCV5	SE	1	21.500	10 31.800	20 42.700	30 52.100	21.0	0.9996	-
ICV1	SE	2	50.600	10 59.100	20 68.000	30 78.800	107.2	0.9984	-
MEW615	SE	1	0.000	10 7.300	20 13.000	30 18.400	14 0.0	0.9973	-
MEW616S	SE	1	13.900	10 18.700	20 25.500	30 33.000	20.5	0.9954	-
CCV4	SE	1	21.100	10 32.000	20 41.100	30 51.900	21.0	0.9994	-
CCV3	SE	1	20.600	10 30.500	20 39.400	30 50.500	20.8	0.9991	-
LCS1	TL	2	47.400	10 59.600	20 68.700	30 81.800	84.6	0.9978	-
CCB3	SE	1	0.000	10 10.600	20 21.700	30 30.900	14 0.0	0.9993	-
LCS	SE	2	50.000	10 61.000	20 71.600	30 79.100	103.6	0.9964	-
MEW621	SE	1	0.000	10 2.300	20 5.400	30 6.500	14 0.0	0.9856	+
ICV	SE	2	48.800	10 58.000	20 67.800	30 76.800	104.0	0.9999	-
CRA	SE	1	1.700	5 6.300	10 11.700	15 17.000	1.4	0.9994	-
CCB1	SE	1	0.000	10 11.600	20 22.100	30 31.900	14 0.0	0.9993	-
CCB2	SE	1	0.000	10 10.300	20 21.600	30 31.800	14 0.0	0.9998	-
CCV2	SE	1	20.500	10 29.800	20 40.100	30 49.400	21.0	0.9998	-
ICB	SE	1	0.000	10 11.100	20 22.100	30 32.100	14 0.0	0.9997	-
CCB2	SE	1	0.000	10 10.000	20 20.200	30 30.400	14 0.0	1.0000	-
ICB	TL	1	0.000	10 9.300	20 20.000	30 28.400	14 0.0	0.9990	-
CCV1	SE	1	20.500	10 30.200	20 40.600	30 49.800	20.9	0.9997	-
ICB1	SE	1	0.000	10 9.900	20 19.900	30 30.700	14 0.0	0.9998	-

U.S. EPA - CLP

8  
STANDARD ADDITION RESULTS

00051

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615

Concentration Units: ug/L

EPA Sample No.	An	Dil	0 ADD ABS	1 ADD CON ABS	2 ADD CON ABS	3 ADD CON ABS	Final Conc.	r	Q
CCV1	TL	1	19.200	10 29.300	20 39.200	30 47.900	20.3	0.9994	-
CCV2	TL	1	19.400	10 30.000	20 39.300	30 48.100	20.8	0.9991	-
CCB1	TL	1	0.000	10 9.900	20 19.500	30 29.400	24 0.0	1.0000	-
CCV2	SE	1	20.200	10 29.800	20 39.400	30 50.700	19.6	0.9992	-
CCB2	TL	1	-1.300	10 9.000	20 19.700	30 29.800	24 0.0	0.9999	-
CCB1	TL	1	-1.400	10 9.300	20 19.800	30 30.700	24 0.0	1.0000	-
CCV1	SE	1	21.000	10 31.600	20 41.000	30 50.300	22.0	0.9995	-
CCB4	SE	1	0.000	10 12.100	20 22.500	30 33.200	14 0.0	0.9994	-
ICB1	TL	1	-1.200	10 9.700	20 20.300	30 30.900	24 0.0	1.0000	-
PB	TL	1	-1.200	10 9.700	20 20.800	30 32.100	24 0.0	1.0000	-
CCB4	TL	1	-1.300	10 9.400	20 19.600	30 29.600	24 0.0	0.9999	-
CCV2	TL	1	18.400	10 26.000	20 37.300	30 47.200	18.0	0.9973	-
ICV	TL	2	47.000	10 56.300	20 67.900	30 78.600	87.4	0.9991	-
CRA	TL	1	0.000	5 8.100	10 14.100	15 19.500	24 0.0	0.9954	-
CCV3	TL	1	19.300	10 30.400	20 38.200	30 50.300	19.3	0.9969	-
CCB3	TL	1	-1.200	10 9.400	20 18.700	30 30.500	24 0.0	0.9990	-
ICV1	TL	2	48.700	10 59.600	20 69.800	30 76.600	105.6	0.9949	+
CCB2	TL	1	0.000	10 8.100	20 17.800	30 27.900	24 0.0	0.9988	-
CCV1	TL	1	19.500	10 30.800	20 39.300	30 51.300	18.9	0.9981	-
CCV4	TL	1	19.000	10 30.900	20 40.500	30 49.200	19.8	0.9974	-
CRA	TL	1	0.000	5 6.100	10 11.600	15 16.800	24 0.0	0.9993	-
MEW618	SE	1	0.000	10 8.400	20 16.300	30 23.200	14 0.0	0.9990	-
MEW619	SE	1	0.000	10 9.800	20 21.000	30 30.800	14 0.0	0.9996	-
MEW621	SE	10	1.500	10 9.800	20 18.900	30 29.700	14 0.0	0.9982	-

00052

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## 8

### STANDARD ADDITION RESULTS

Lab Name: ROCKY MOUNTAIN ANALYTICAL

Contract: 68-01-7476

**Lab Code: ENSECO**

**Case No.: 9288**

**SAS No.: 3554E**

**SDG No.: MEW615**

**Concentration Units: ug/L**

00053

U.S. EPA - CLP

9  
ICP SERIAL DILUTIONS

EPA SAMPLE NO.

Lab Name: ROCKY MOUNTAIN ANALYTICAL Contract: 68-01-7476MEW623LLab Code: ENSECO Case No.: 9288 SAS No.: 3554E SDG No.: MEW615Matrix (soil/water): WATERLevel (low/med): LOW

Concentration Units: ug/L

Analyte	Initial Sample Result (I)	C	Serial Dilution Result (S)	C	% Difference	Q	M
Aluminum	216	-	187	B	13.4	-	NR
Antimony	-	-	-	-	-	-	-
Arsenic	-	-	-	-	-	-	-
Barium	90.6	-	86.6	B	4.4	-	NR
Beryllium	2.0	U	10.0	U	-	-	NR
Cadmium	-	-	-	-	-	-	-
Calcium	85470	-	85760	-	0.3	-	P
Chromium	4.0	U	20.0	U	-	-	NR
Cobalt	120	U	113	U	5.8	-	NR
Copper	10.0	U	50.0	U	-	-	NR
Iron	76780	-	77360	-	0.8	-	P
Lead	-	-	-	-	-	-	-
Magnesium	26990	-	26860	-	0.5	-	P
Manganese	517	-	501	-	3.1	-	P
Mercury	-	-	-	-	-	-	-
Nickel	342	-	327	B	4.4	-	NR
Potassium	1164	B	1115	B	4.2	-	NR
Selenium	-	-	-	-	-	-	-
Silver	4.0	U	20.0	U	-	-	NR
Sodium	3600	-	6800	U	100	-	NR
Thallium	-	-	-	-	-	-	-
Vanadium	4.0	U	20.0	U	-	-	NR
Zinc	477	-	486	-	1.9	-	P

00054

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10  
HOLDING TIMES

**Lab Name:** ROCKY MOUNTAIN ANALYTICAL

**Contract:** 68-01-7476

**Lab Code:** ENSECO

**Case No.: 9288**

SAS No.: 3554E

**SDG No.:** MEW615

11  
INSTRUMENT DETECTION LIMITS (QUARTERLY)

00055

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9188SAS No.: 3554ESDG No.: new 615ICP ID Number: JIA9000Date: 01/26/88

Flame AA ID Number: \_\_\_\_\_

Furnace AA ID Number: \_\_\_\_\_

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum	237.31		100	25	P
Antimony			5		
Arsenic			5		
Barium	233.53		50	2.0	P
Beryllium	313.04		5	2.0	P
Cadmium			0.5		
Calcium	370.60		1000	81	P
Chromium	267.72		10	4.0	P
Cobalt	228.62		10	4.0	P
Copper	324.75		10	10	P
Iron	259.84		100	14	P
Lead	220.35		2	17	P
Magnesium	279.08		1000	81	P
Manganese	294.92		10	6.0	P
Mercury			0.2		
Nickel	231.60		20	10	P
Potassium	766.49		2000	95	P
Selenium			2		
Silver	328.07		5	4.0	P
Sodium	330.24		1000	1360	P
Thallium			2		
Vanadium	292.40		10	4.0	P
Zinc	213.80		20	3.0	P

## Comments:

SPECTRO-PRODUCTS HG-3 SPECTROPHOTOMETER USED FOR MANUAL COLD VAPOR DETERMINATION

11  
INSTRUMENT DETECTION LIMITS (QUARTERLY)

00056

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554E SDG No.: NEW615

ICP ID Number:

JA9000Date: 01/26/88

Flame AA ID Number:

Furnace AA ID Number: PE2380

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			100		
Antimony			5		
Arsenic	193.70	BD	5	3.0	F
Barium			50		
Beryllium			5		
Cadmium	228.80	BD	0.5	0.1	F
Calcium			1000		
Chromium			10		
Cobalt			10		
Copper			10		
Iron			100		
Lead	283.30	BD	2	1.0	F
Magnesium			1000		
Manganese			10		
Mercury	253.70	BD	0.2	0.1	CV
Nickel			20		
Potassium			2000		
Selenium	196.03	BD	2	2.0	F
Silver			5		
Sodium			1000		
Thallium	276.80	BD	2	2.0	F
Vanadium			10		
Zinc			20		

## Comments:

SPECTRO-PRODUCTS HG-3 SPECTROPHOTOMETER USED FOR MANUAL COLD VAPOR DETERMINATION

11  
INSTRUMENT DETECTION LIMITS (QUARTERLY)

Lab Name: ROCKY MOUNTAIN ANALYTICALContract: 68-01-7476Lab Code: ENSECOCase No.: 9288SAS No.: 3554ESDG No.: MEW615

ICP ID Number:

JA9000Date: 01/26/88

Flame AA ID Number:

Furnace AA ID Number: 25000

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			100		
Antimony	217.80	BZ	5	2.0	F
Arsenic	193.70	BZ	5	1.0	F
Barium			50		
Beryllium			5		
Cadmium			0.5		
Calcium			1000		
Chromium			10		
Cobalt			10		
Copper			10		
Iron			100		
Lead			2		
Magnesium			1000		
Manganese			10		
Mercury			0.2		
Nickel			20		
Potassium			2000		
Selenium	196.03	BZ	2	1.0	F
Silver			5		
Sodium			1000		
Thallium			2		
Vanadium			10		
Zinc			20		

## Comments:

SPECTRO-PRODUCTS HG-3 SPECTROPHOTOMETER USED FOR MANUAL COLD VAPOR DETERMINATION

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In Reference to Case No(s):

SAS 3727E

Contract Laboratory Program  
REGIONAL/LABORATORY COMMUNICATION SYSTEM

QC 1

Telephone Record Log

Date of Call: 5.11.88.

Laboratory Name: PBS & J

Lab Contact: Kimberly Leenihis

Region: V

Regional Contact: Addie Leenihis

Call Initiated By: Laboratory  Region

In reference to data for the following sample number(s):

E01-E09

Summary of Questions/Issues Discussed:

1. Acidity - why you did not add 5ml 0.2NH<sub>2</sub>SO<sub>4</sub> to the sample (method 305.1)
2. Alkalinity - what was your lab. blank volume (end) 25mils or 50mils
3. How did you convert short units for chloride, NO<sub>3</sub>-NO<sub>2</sub>; sulfate

Summary of Resolution:

1. Lab found problems with add. no 5ml 11.5c
2. will be corrected and sent to the Regio
3. Producer's copy was sent

Signature

5.12.88.  
Date

Distribution: (1) Lab Copy, (2) Region Copy, (3) SMO Copy

SF 5011

In Reference to Case No(s):

SAS 3727E

Contract Laboratory Program  
REGIONAL/LABORATORY COMMUNICATION SYSTEM  
Telephone Record Log

Date of Call: 4/5/88

Laboratory Name: PBS&J

Lab Contact: Tom French

Region: IV

Regional Contact: Jeb Livingston, Tony Warky (SMO)

Call Initiated By:  Laboratory  Region

In reference to data for the following sample number(s):

3727E-C

Summary of Questions/Issues Discussed:

Received eight containers from San C. City Water and Sewer Board. How do we treat effluent? 12 to 14 ppm chlorine. Extra chlorine at 12-15 ppm for 20 minutes.  
But . . . a chlorine emergency situation  
but . . . i.e., if SMO's

Summary of Resolution:

Treat sample as an individual sample, use extra chlorine until 14 ppm.

Signature

French

Date

4/6/88

Distribution: (1) Lab Copy, (2) Region Copy, (3) SMO Copy

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

DATE: Jay Thacker 5/13/88

BJECT: Review of Region V CLP Data  
Received for Review on 5-2-88

FROM: Curtis Ross ,Director (5SCRL)  
Central Regional Laboratory Jay Thacker

TO: Data User : MPCA  
DAVE KOVBSKY

We have reviewed the data for the following case(s).

SITE NAME: HADER GROUND WATER CONT. SMO case No. SAS3727  
No.of D.U/Activity  
EPA Data Set No. SF5011 Samples: 9 Numbers Y905/C722ZZ

CRL No. 88YL06S01-S08,D05

SMO Traffic No. E01-E09 Hrs.required  
CLP Laboratory: PBS&J for Review 3

Following are our findings:

This review covers nine low water samples analyzed for sulfate, chloride, nitrate/nitrite, alkalinity, acidity and pH.

All QC data are acceptable.

John Levin

5.12.88

- (X) Data are acceptable for use.  
( ) Data are acceptable for use with qualifications noted above.  
( ) Data are preliminary-pending verification by Contract Laboratory.  
( ) Data are unacceptable.

cc: Carla Dempsey, Quality Assurance Officer, EPA Support Services  
James Petty, Chief Quality Assurance Research, EMSL, Las Vegas

U.S.A EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

A large, faint stamp is visible in the background. It contains the word "RECEIVED" at the top, followed by the date "MAY 02 1988" in the center, and "U.S. EPA CENTRAL REGIONAL LAB" at the bottom.

Date 4/26/88

COVER PAGE

Lab Name Post, Buckley, Schuh & Jernigan

Case No. SAS 3727E

SOW No. 7/85

Q.C. Report No. 1

### Sample Numbers

EPA No.	Lab ID No.	EPA No.	Lab Id No.
3727E-01	8804014-01	3727E-09	8804014-09
3727E-02	8804014-02		
3727E-03	8804014-03		
3727E-04	8804014-04		
3727E-05	8804014-05		
3727E-06	8804014-06		
3727E-07	8804014-07		
3727E-08	8804014-08		
Comments:			

ICP interelement and background corrections applied? Yes  No   
If yes, corrections applied before  or after  generation of raw data.

### **Footnotes:**

NR - Not required by contract at this time

**Form I:**

**Value** - If the result is a value greater than or equal to the instrument detection limit but less than the contract-required detection limit, report the value in brackets (i.e., [10]). Indicate the analytical method used with P (for ICP), A (for Flame AA) or F (for Furnace AA).

U - Indicates element was analyzed for but not detected. Report with the instrument detection limit value (e.g., 100).

E - Indicates a value estimated or not reported due to the presence of interference. Explanatory note included on cover page.

s - Indicates value determined by Method of Standard Addition.

N - Indicates spike sample recovery is not within control limits.

\* - Indicates duplicate analysis is not within control limits.

+ - Indicates the correlation coefficient for method of standardization.

M - Indicates duplicate injection results exceeded control limits.

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No. |

| 3727E-01 |

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-01

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium \_\_\_\_\_

Matrix: Water  Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L) <sup>ppm</sup> ug/g or mg/kg dry weight (Circle One)

1. Sulfate 605 P 10. \_\_\_\_\_

2. Nitrate & Nitrite 0.10 P 11. \_\_\_\_\_

3. Chloride 0.50 F 12. \_\_\_\_\_

4. Alkalinity 122 P 13. \_\_\_\_\_

5. Acidity \_\_\_\_\_ P 14. \_\_\_\_\_

6. PH 6.4 P 15. \_\_\_\_\_

7. \_\_\_\_\_ P 16. \_\_\_\_\_

8. \_\_\_\_\_ P 17. \_\_\_\_\_

9. \_\_\_\_\_ Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless clear ( $\text{NO}_3 + \text{NO}_2$ )  
Orange, Cloudy (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager

K.A. Kuehne

IFB Amend One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.  
3727E-02

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-02

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium

Matrix: Water  Soil  Sludge  Other

(mg/l or ug/g) or mg/kg dry weight (Circle One)

1. Sulfate	<u>310</u>	P	10.	
2. Nitrate & Nitrite	<u>0.10</u>	P	11.	
3. Chloride	<u>0.50</u>	F	12.	
4. Alkalinity	<u>231</u>	P	13.	
5. Acidity		P	14.	
6. PH	<u>6.8</u>	P	15.	
7.		P	16.	
8.		P	17.	
9.			Percent Solids (%)	

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless, clear N<sub>2</sub>O<sub>3</sub>+N<sub>2</sub>O<sub>5</sub>  
Yellow, clear (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager K. J. Schuh

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-577-2490

EPA Sample No.  
3727E-03

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan CASE NO. SAS 3727E

SOW NO. 7/85 Lab Receipt Date 4/2/88

LAB SAMPLE ID. NO. 8804014-03 Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low X Medium \_\_\_\_\_

Matrix: Water X Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L) 41250 mg/kg dry weight or mg/kg dry weight (Circle One)

- |                      |             |           |   |     |                          |
|----------------------|-------------|-----------|---|-----|--------------------------|
| 1. Sulfate           | <u>413</u>  | <u>19</u> | P | 10. |                          |
| 2. Nitrate & Nitrite | <u>0.10</u> |           | P | 11. |                          |
| 3. Chloride          | <u>0.50</u> |           | F | 12. |                          |
| 4. Alkalinity        | <u>277</u>  |           | P | 13. |                          |
| 5. Acidity           |             |           | P | 14. |                          |
| 6. PH                | <u>7.2</u>  |           | P | 15. |                          |
| 7.                   |             |           | P | 16. |                          |
| 8.                   |             |           | P | 17. |                          |
| 9.                   |             |           |   |     | Percent Solids (%) _____ |

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless, clear (N<sub>2</sub>O<sub>3</sub>+N<sub>2</sub>O<sub>2</sub>)  
White, clear (Alkalinity, Acidity PH, Sulfate, Chloride)

Lab Manager J.A. Kunkel

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-577-2490

EPA Sample No.  
3727E-04

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-04

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low X Medium \_\_\_\_\_

Matrix: Water X Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/l or ug/g) or mg/kg dry weight (Circle One)

- |                                 |             |   |                          |       |
|---------------------------------|-------------|---|--------------------------|-------|
| 1. <u>Sulfate</u>               | <u>315</u>  | P | 10.                      | _____ |
| 2. <u>Nitrate &amp; Nitrite</u> | <u>0.10</u> | P | 11.                      | _____ |
| 3. <u>Chloride</u>              | <u>0.50</u> | F | 12.                      | _____ |
| 4. <u>Alkalinity</u>            | <u>127</u>  | P | 13.                      | _____ |
| 5. <u>Acidity</u>               |             | P | 14.                      | _____ |
| 6. <u>pH</u>                    | <u>6.3</u>  | P | 15.                      | _____ |
| 7.                              |             | P | 16.                      | _____ |
| 8.                              |             | P | 17.                      | _____ |
| 9.                              |             |   | Percent Solids (%) _____ |       |

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless, clear (NO<sub>3</sub>+NO<sub>2</sub>)  
Orange, cloudy (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager K. A. Kunkle

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-577-2490

EPA Sample No.  
3727E-05

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-05

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low X Medium \_\_\_\_\_

Matrix: Water X Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L <sup>w/w</sup>) or mg/kg dry weight (Circle One)

- |                      |             |   |                          |       |
|----------------------|-------------|---|--------------------------|-------|
| 1. Sulfate           | <u>136</u>  | P | 10.                      | _____ |
| 2. Nitrate & Nitrite | <u>0.10</u> | P | 11.                      | _____ |
| 3. Chloride          | <u>0.50</u> | F | 12.                      | _____ |
| 4. Alkalinity        | <u>324</u>  | P | 13.                      | _____ |
| 5. Acidity           |             | P | 14.                      | _____ |
| 6. PH                | <u>7.3</u>  | P | 15.                      | _____ |
| 7.                   |             | P | 16.                      | _____ |
| 8.                   |             | P | 17.                      | _____ |
| 9.                   |             |   | Percent Solids (%) _____ |       |

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments Colorless, clear (NO<sub>3</sub>+NO<sub>2</sub>)  
Orange, clear (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager

T. A. Kunkel

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.

3727E-06

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/2/88

LAB SAMPLE ID. NO. 8804014-06

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium \_\_\_\_\_

Matrix: Water  Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L <sup>wet weight</sup> or mg/kg dry weight) (Circle One)

1. Sulfate	890	P	10.	
2. Nitrate & Nitrite	0.10	P	11.	
3. Chloride	0.50	F	12.	
4. Alkalinity	2.00	P	13.	
5. Acidity	576	P	14.	
6. PH	4.6	P	15.	
7.		P	16.	
8.		P	17.	
9.				Percent Solids (%)

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however

Comments: Colorless, clear (No ENs)  
White, cloudy (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager

Kelli Miller

JFB Amend One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-577-2490

EPA Sample No.  
3727E-07

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-07

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low X Medium \_\_\_\_\_

Matrix: Water X Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/l) <sup>41581 mg</sup> or mg/kg dry weight (Circle One)

- |                      |             |                                |                          |       |       |
|----------------------|-------------|--------------------------------|--------------------------|-------|-------|
| 1. Sulfate           | <u>24</u>   | P                              | 10.                      | _____ |       |
| 2. Nitrate & Nitrite | <u>0.10</u> | P                              | 11.                      | _____ |       |
| 3. Chloride          | <u>3.5</u>  | <sup>41581 mg</sup> <u>3.4</u> | F                        | 12.   | _____ |
| 4. Alkalinity        | <u>340</u>  | P                              | 13.                      | _____ |       |
| 5. Acidity           |             | P                              | 14.                      | _____ |       |
| 6. PH                | <u>7.0</u>  | P                              | 15.                      | _____ |       |
| 7.                   |             | P                              | 16.                      | _____ |       |
| 8.                   |             | P                              | 17.                      | _____ |       |
| 9.                   |             |                                | Percent Solids (%) _____ |       |       |

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: (Colorless, clear (N<sub>2</sub>+N<sub>2</sub>O))  
Yellow, clear (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager J.T. Kuehne

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.

3727E-08

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/2/88

LAB SAMPLE ID. NO. 8804014-08

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium \_\_\_\_\_

Matrix: Water  Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/l) <sup>ppm</sup> or mg/kg dry weight (Circle One)

- |                      |      |   |                          |  |
|----------------------|------|---|--------------------------|--|
| 1. Sulfate           | 314  | P | 10.                      |  |
| 2. Nitrate & Nitrite | 0.10 | P | 11.                      |  |
| 3. Chloride          | 1.2  | F | 12.                      |  |
| 4. Alkalinity        | 134  | P | 13.                      |  |
| 5. Acidity           |      | P | 14.                      |  |
| 6. PH                | 6.5  | P | 15.                      |  |
| 7.                   |      | P | 16.                      |  |
| 8.                   |      | P | 17.                      |  |
| 9.                   |      |   | Percent Solids (%) _____ |  |

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless, clear ( $\text{NO}_3 + \text{NO}_2$ )  
Orange, clear (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager X. J. Finch

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No. |

| 3727E-09 |

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/2/88

LAB SAMPLE ID. NO. 8804014-09

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium \_\_\_\_\_

Matrix: Water  Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L ug/g or mg/kg dry weight (Circle One))

- |                      |      |   |                          |       |
|----------------------|------|---|--------------------------|-------|
| 1. Sulfate           | 139  | P | 10.                      | _____ |
| 2. Nitrate & Nitrite | 0.10 | P | 11.                      | _____ |
| 3. Chloride          | 0.50 | F | 12.                      | _____ |
| 4. Alkalinity        | 329  | P | 13.                      | _____ |
| 5. Acidity           |      | P | 14.                      | _____ |
| 6. PH                | 7.3  | P | 15.                      | _____ |
| 7.                   |      | P | 16.                      | _____ |
| 8.                   |      | P | 17.                      | _____ |
| 9.                   |      |   | Percent Solids (%) _____ |       |

Footnotes For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments Colored, clear (N<sub>2</sub>O<sub>3</sub>+N<sub>2</sub>) Orange, clear (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager JGJ/Jernigan

IFB Amend. One

Form III

Q.C. Report No. 1

BLANKS

LAB NAME Post, Buckley, Schuh & Jernigan  
DATE 4/26/88

CASE NO. SAS 3727E  
UNITS ug/l mg/L

Compound	Initial Calibration Blank Value	Continuing Calibration				Prep. Blank	
		1	2	3	4	Matrix	Matrix
						1	2
1. Sulfate	5.0u	5.0u					
2. NO <sub>3</sub> - NO <sub>2</sub>	0.1u	0.1u				0.1u	0.1u
3. Chloride	0.5u	0.5u					
4. Alkalinity	10u	10u					
5. Acidity							
6. PH							
7.							
8.							
9.							
10.							
11.							
12.							
13.							
14.							
15.							
16.							
17.							
18.							
19.							
20.							
21.							
22.							
23.							
Other:							
Cyanide							

<sup>1</sup> Reporting Units: aqueous, ug/l; solid, mg/kg

Form V *4/26/88*

Q.C. Report No. 1

SPIKE SAMPLE RECOVERY

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

DATE

4/26/88

EPA Sample No. 3727E-01

Lab Sample ID No. 8804014-01

Units: <sup>mg/l</sup> mg/kg mg/L

Matrix LA

Compound	Control Limit %R	Spiked Sample Result (SSR)	Sample Result(SR)	Spiked Added (SA)	%R
1. Sulfate	85-115	1450	605	800	106
2. N03 - NO2	"	0.8	0.10	0.8	100
3. Chloride	"	4.5	0.50	4.0	112
4. Alkalinity	"	NR			
5. Acidity	"				
6. PH	"				
7.	"				
8.	"				
9.	"				
10.	"				
11.	"				
12.	"				
13.	"				
14.	"				
15.	"				
16.	"				
17.	"				
18.	"				
19.	"				
20.	"				
21.	"				
22.	"				
23.	"				
Other					
Cyanide					

<sup>1</sup> %R = [ ( SSR - SR ) / SA ] x 100

" N " - out of control

' NR "-Not required

Comments.

Form V b

Q.C. Report No. 1

DUPLICATES

LAB NAME Post, Buckley, Schuh & Jernigan  
DATE 4/26/88

CASE NO. SAS 3727E  
EPA Sample No. 3727E-06  
Lab Sample ID No. 8804014-06  
Units: mg/L

Matrix LA

Compound	Control Limit1	Sample (S)	Duplicate (D)	RPD2
1. Sulfate				
2. NO <sub>3</sub> - NO <sub>2</sub>				
3. Chloride				
4. Alkalinity				
5. Acidity		576	578	0.3
6. PH		4.6	4.6	0.0
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
Other:				
Cyanide				

\* Out of Control

<sup>1</sup> To be added at a later date

<sup>2</sup> RPD = [ S - D / ((S + D)/2) ] x 100

NC - Non calculable RPD due to value(s) less than CRDL

Form M C

Q.C. Report No. 1

DUPLICATES

LAB NAME Post, Buckley, Schuh & Jernigan  
DATE 4/26/88

CASE NO. SAS 3727E  
EPA Sample No. 3727E-09  
Lab Sample ID No. 8804014-09  
Units: mg/L

Matrix LA

Compound	Control Limit1	Sample (S)	Duplicate (D)	RPD2
1. Sulfate				
2. NO <sub>3</sub> - NO <sub>2</sub>				
3. Chloride				
4. Alkalinity	+10%	329	328	0.3
5. Acidity				
6. PH				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
Other:				
Cyanide				

\* Out of Control

<sup>1</sup> To be added at a later date.

$$^2 \text{ RPD} = [ (S - D) / ((S + D)/2) ] \times 100$$

NC - Non calculable RPD due to value(s) less than CRDL

Form V d

Q.C. Report No. 1

DUPLICATES

LAB NAME Post, Buckley, Schuh & Jernigan  
DATE 4/26/88

CASE NO. SAS 3727E  
EPA Sample No. 3727E-01  
Lab Sample ID No. 8804014-01  
Units: mg/l

Matrix LA

Compound	Control Limit <sup>1</sup>	Sample (S)	Duplicate (D)	RPD <sup>2</sup>
1. Sulfate	<u>±10%</u>	<u>588</u>	<u>594</u>	<u>1.0</u>
2. NO <sub>3</sub> - NO <sub>2</sub>		<u>0.1u</u>	<u>0.1u</u>	<u>NC</u>
3. Chloride		<u>0.5u</u>	<u>0.5u</u>	<u>NC</u>
4. Alkalinity				
5. Acidity				
6. PH				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
Other:				
Cyanide				

\* Out of Control

<sup>1</sup> To be added at a later date.

<sup>2</sup> RPD = [ (S - D) / ((S + D)/2) ] x 100

NC - Non calculable RPD due to value(s) less than CRDL

Form VII  
 Q.C. Report No. 1  
 INSTRUMENT DETECTION LIMITS AND  
 LABORATORY CONTROL SAMPLE  
 LAB NAME Post, Buckley, Schuh & Jernigan CASE NO. SAS 3727E DATE 4/26/86

LCS NO. SAS 3727E

Metals: Compound	Required Detection Limits (CRDL)- ug/l		Instrument Detection Limits (IDL)- ug/l		Lab Control Sample mg/kg (circle one)	
	Mg/L	ID#	ICP/AA	Furnace	ID#	True Found %R
Metals:						
1. Sulfate	5.0		5.0		20	20 100
2. NO <sub>3</sub> - NO <sub>2</sub>	0.1		0.10		2.0	2.0 100
3. Chloride	0.5		0.5		52	52 100
4. Alkalinity	10		10		25	25 100
5. Acidity	± 10.		± 10		6.0	6.0 100
6. PH	0.1		0.1		NR	6.0 6.0 100
7.						
8.						
9.						
10.						
11.						
12.						
13.						
14.						
15.						
16.						
17.						
18.						
19.						
20.						
21.						
22.						
23.						
Other						
Cyanide	10		NR	NR		

NR - Not required

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

8 Solid Waste Div.  
MPCA Ground Water

DATE: Jay Thakker 5/13/88 MW 19.8

SUBJECT: Review of Region V CLP Data  
Received for Review on 5-2-88

FROM: Curtis Ross ,Director (SSCRL)  
Central Regional Laboratory Jay Thakker

TO: Data User : MPCA  
DAVE KOVBSKY

REVIEWED

We have reviewed the data for the following case(s).

SITE NAME: HADER GROUND WATER CONT. SMO case No. SAS3727

EPA Data Set No. SF5011 No.of Samples: 9 D.U/Activity Numbers Y905/C722ZZ

CRL No. 88YL06S01-S08,D05

SMO Traffic No. E01-E09

CLP Laboratory: PBS&J Hrs.required for Review 3

Following are our findings:

This review covers nine low water samples analyzed for sulfate, chloride, nitrate/nitrite, alkalinity, acidity and pH.

All QC data are acceptable.

John Levin

5.12.88

- Data are acceptable for use.
- Data are acceptable for use with qualifications noted above.
- Data are preliminary-pending verification by Contract Laboratory.
- Data are unacceptable.

cc: Carla Dempsey, Quality Assurance Officer, EPA Support Services  
James Petty, Chief Quality Assurance Research, EMSL, Las Vegas

*U.S.A. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490*

*U.S. EPA CENTRAL  
REGIONAL LAB*

*Date 4/26/88*

COVER PAGE  
INORGANIC ANALYSES DATA PACKAGE

Lab Name Post, Buckley, Schuh & Jernigan

Case No. SAS 3727E

SOW No. 7/85

Q.C. Report No. 1

Sample Numbers

EPA No.	Lab ID No.	EPA No.	Lab Id No.
3727E-01	8804014-01	3727E-09	8804014-09
3727E-02	8804014-02		
3727E-03	8804014-03		
3727E-04	8804014-04		
3727E-05	8804014-05		
3727E-06	8804014-06		
3727E-07	8804014-07		
3727E-08	8804014-08		

Comments:

ICP interelement and background corrections applied? Yes  No   
If yes, corrections applied before  or after  generation of raw data.

Footnotes:

NR - Not required by contract at this time

Form I:

Value - If the result is a value greater than or equal to the instrument detection limit but less than the contract-required detection limit, report the value in brackets (i.e., [10]). Indicate the analytical method used with P (for ICP), A (for Flame AA) or F (for Furnace AA).

U - Indicates element was analyzed for but not detected. Report with the instrument detection limit value (e.g., 10U).

E - Indicates a value estimated or not reported due to the presence of interference. Explanatory note included on cover page.

S - Indicates value determined by Method of Standard Addition.

N - Indicates spike sample recovery is not within control limits.

\* - Indicates duplicate analysis is not within control limits.

+ - Indicates the correlation coefficient for method of standard addition is less than 0.995.

M - Indicates duplicate injection results exceeded control limits.

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.  
3727E-01

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-01

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low X Medium \_\_\_\_\_

Matrix: Water X Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L)ug/g or mg/kg dry weight (Circle One)

1. Sulfate 605 P 10. \_\_\_\_\_

2. Nitrate & Nitrite 0.10 P 11. \_\_\_\_\_

3. Chloride 0.50 F 12. \_\_\_\_\_

4. Alkalinity 122 P 13. \_\_\_\_\_

5. Acidity \_\_\_\_\_ P 14. \_\_\_\_\_

6. PH 6.4 P 15. \_\_\_\_\_

7. \_\_\_\_\_ P 16. \_\_\_\_\_

8. \_\_\_\_\_ P 17. \_\_\_\_\_

9. \_\_\_\_\_ Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless, clear (NO<sub>3</sub>+NO<sub>2</sub>)  
Orange, cloudy (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager K.A.Kimball

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

|EPA Sample No.  
| 3727E-02 |

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post: Buckley, Schuh & Jernigan CASE NO SAS 3727E

SOW NO. 7/85 Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-02 Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium

Matrix: Water  Soil  Sludge  Other

(mg/l wet wt or mg/kg dry weight (Circle One))

1. Sulfate	310	P	10.	
2. Nitrate & Nitrite	0.10	P	11.	
3. Chloride	0.50	F	12.	
4. Alkalinity	231	P	13.	
5. Acidity		P	14.	
6. PH	6.8	P	15.	
7.		P	16.	
8.		P	17.	
9.			Percent Solids (%)	

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless, clear ( $\text{NO}_3 + \text{N}_2$ )  
Yellow, clear (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager K.T. Jernigan

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.  
3727E-03

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan CASE NO SAS 3727E

SOW NO. 7/85 Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-03 Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium

Matrix: Water  Soil  Sludge  Other

(mg/L) 412546 μm³ or mg/kg dry weight (Circle One)

1. Sulfate 313 19 P 10. \_\_\_\_\_

2. Nitrate & Nitrite 0.10 P 11. \_\_\_\_\_

3. Chloride 0.50 F 12. \_\_\_\_\_

4. Alkalinity 277 P 13. \_\_\_\_\_

5. Acidity \_\_\_\_\_ P 14. \_\_\_\_\_

6. PH 7.2 P 15. \_\_\_\_\_

7. \_\_\_\_\_ P 16. \_\_\_\_\_

8. \_\_\_\_\_ P 17. \_\_\_\_\_

9. \_\_\_\_\_ Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colorless, clear ( $\text{Na}_2\text{O} + \text{Na}_2\text{S}_2\text{O}_3$ )  
White, clear (Alkalinity, Acidity pH, Sulfate, Chloride)

Lab Manager

P. A. Kunkle

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-577-2490

EPA Sample No.  
3727E-04

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO SAS 3727E

SOW NO 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-04

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium \_\_\_\_\_

Matrix: Water  Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(<sup>ug/L</sup> or mg/kg dry weight) (Circle One)

1. Sulfate 315 P 10. \_\_\_\_\_

2. Nitrate & Nitrite 0.10 P 11. \_\_\_\_\_

3. Chloride 0.50 F 12. \_\_\_\_\_

4. Alkalinity 127 P 13. \_\_\_\_\_

5. Acidity \_\_\_\_\_ P 14. \_\_\_\_\_

6. PH 6.3 P 15. \_\_\_\_\_

7. \_\_\_\_\_ P 16. \_\_\_\_\_

8. \_\_\_\_\_ P 17. \_\_\_\_\_

9. \_\_\_\_\_ Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments. Colorless, clear (NO<sub>3</sub>+NO<sub>2</sub>)  
Orange, cloudy (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager

K. Kunkel

IFB Amend One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.  
3727E-05

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-05

Q.C REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium \_\_\_\_\_

Matrix: Water  Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L wet weight or mg/kg dry weight) (Circle One)

- |                      |             |   |     |                          |
|----------------------|-------------|---|-----|--------------------------|
| 1. Sulfate           | <u>136</u>  | P | 10. | _____                    |
| 2. Nitrate & Nitrite | <u>0.10</u> | P | 11. | _____                    |
| 3. Chloride          | <u>0.50</u> | F | 12. | _____                    |
| 4. Alkalinity        | <u>324</u>  | P | 13. | _____                    |
| 5. Acidity           |             | P | 14. | _____                    |
| 6. PH                | <u>7.3</u>  | P | 15. | _____                    |
| 7.                   |             | P | 16. | _____                    |
| 8.                   |             | P | 17. | _____                    |
| 9.                   |             |   |     | Percent Solids (%) _____ |

Footnotes. For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments. Colorless, clear (No<sub>3</sub>+No<sub>2</sub>)  
Orange, clear (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager K. Kinkade

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.  
3727E-06

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO. 8804014-06

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium

Matrix: Water  Soil  Sludge  Other

(<sup>mg/l</sup> or <sup>ug/g</sup>) or mg/kg dry weight (Circle One)

- |                      |             |   |                    |  |
|----------------------|-------------|---|--------------------|--|
| 1. Sulfate           | <u>890</u>  | P | 10.                |  |
| 2. Nitrate & Nitrite | <u>0.1</u>  | P | 11.                |  |
| 3. Chloride          | <u>0.50</u> | F | 12.                |  |
| 4. Alkalinity        | <u>2.00</u> | P | 13.                |  |
| 5. Acidity           | <u>576</u>  | P | 14.                |  |
| 6. PH                | <u>4.6</u>  | P | 15.                |  |
| 7.                   |             | P | 16.                |  |
| 8.                   |             | P | 17.                |  |
| 9.                   |             |   | Percent Solids (%) |  |

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Colored, clear ( $\text{NO}_3 + \text{NH}_3$ )  
White, cloudy (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager K. J. Miller

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.  
3727E-07

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/21/88

LAB SAMPLE ID. NO 8804014-07

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low X Medium \_\_\_\_\_

Matrix: Water X Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L or mg/kg dry weight) (Circle One)

- |                      |              |   |                    |  |
|----------------------|--------------|---|--------------------|--|
| 1. Sulfate           | <u>24</u>    | P | 10.                |  |
| 2. Nitrate & Nitrite | <u>0.10</u>  | P | 11.                |  |
| 3. Chloride          | <u>353.4</u> | F | 12.                |  |
| 4. Alkalinity        | <u>340</u>   | P | 13.                |  |
| 5. Acidity           |              | P | 14.                |  |
| 6. PH                | <u>7.0</u>   | P | 15.                |  |
| 7.                   |              | P | 16.                |  |
| 8.                   |              | P | 17.                |  |
| 9.                   |              |   | Percent Solids (%) |  |

Footnotes For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments. Colorless, clear (N<sub>2</sub>O + N<sub>2</sub>)  
Yellow, clear (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager

JGKunluc

IFB Amend. One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No. 1

13727E-08

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

SOW NO. 7/85

Lab Receipt Date 4/2/88

LAB SAMPLE ID. NO. 8804014-08

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration: Low  Medium \_\_\_\_\_

Matrix: Water  Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/l, ug/g, or mg/kg dry weight) (Circle One)

1. Sulfate	314	P	10.	
2. Nitrate & Nitrite	0.10	P	11.	
3. Chloride	1.2	F	12.	
4. Alkalinity	134	P	13.	
5. Acidity		P	14.	
6. PH	6.5	P	15.	
7.		P	16.	
8.		P	17.	
9.				Percent Solids (%)

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however

Comments: Colored, clear ( $\text{NO}_3 + \text{NO}_2$ )  
Orange, clear (Alkalinity, Acidity, PH, Sulfate, Chloride)

Lab Manager

IFB Amend One

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 PTS: 8-577-2490

EPA Sample No.  
3727E-09

Date 4/26/88

INORGANIC ANALYSIS DATA SHEET

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO SAS 3727E

SOW NO 7/85

Lab Receipt Date 4/2/88

LAB SAMPLE ID NO 8804014-09

Q.C. REPORT NO. 1

Elements Identified and Measured

Concentration. Low X Medium \_\_\_\_\_

Matrix: Water X Soil \_\_\_\_\_ Sludge \_\_\_\_\_ Other \_\_\_\_\_

(mg/L ~~ug/g~~ or mg/kg dry weight) (Circle One)

- |                      |             |   |                    |       |
|----------------------|-------------|---|--------------------|-------|
| 1. Sulfate           | <u>139</u>  | P | 10.                | _____ |
| 2. Nitrate & Nitrite | <u>0.10</u> | P | 11.                | _____ |
| 3. Chloride          | <u>0.50</u> | F | 12.                | _____ |
| 4. Alkalinity        | <u>329</u>  | P | 13.                | _____ |
| 5. Acidity           |             | P | 14.                | _____ |
| 6. PH                | <u>7.3</u>  | P | 15.                | _____ |
| 7.                   |             | P | 16.                | _____ |
| 8.                   |             | P | 17.                | _____ |
| 9.                   |             |   | Percent Solids (%) | _____ |

Footnotes For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments Colored, clear (N<sub>2</sub>O<sub>3</sub>+N<sub>2</sub>) Orange, clear (Alkalinity, Acidity, pH, Sulfate, Chloride)

Lab Manager JG Jernigan

IFB Amend. One

Form III

Q.C. Report No 1

BLANKS

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO S.A.S. 3727 E

DATE 4/26/88

UNITS ug/ml

Compound	Initial Calibration Blank Value	Continuing Calibration				Prep. Blank Matrix Matrix
		1	2	3	4	
1. Sulfate	5.06	5.06				
2. NO <sub>3</sub> - NO <sub>2</sub>	0.10	0.10				0.10 0.10
3. Chloride	0.50	0.50				
4. Alkalinity	100	100				
5. Acidity						
6. PH						
7.						
8.						
9.						
10.						
11.						
12.						
13.						
14.						
15.						
16.						
17.						
18.						
19.						
20.						
21.						
22.						
23.						
Other:						
Cyanide						

<sup>1</sup> Reporting Units: aqueous, ug/L; solid, mg/kg

Form V *d4L*

Q.C. Report No. 1

SPIKE SAMPLE RECOVERY

LAB NAME Post, Buckley, Schuh & Jernigan

CASE NO. SAS 3727E

DATE 4/26/88

EPA Sample No. 3727E-01

Lab Sample ID No. 8804014-01

Units: <sup>#</sup> ~~kg/L~~ mg/L

Matrix LA

Compound	Control Limit *R	Spiked Sample Result (SSR)	Sample Result(SR)	Spiked Added (SA)	%R
1. Sulfate	85-115	1450	605	800	106
2. NO <sub>3</sub> - NO <sub>2</sub>	"	0.8	0.10	0.8	100
3. Chloride	"	4.5	0.50	4.0	112
4. Alkalinity	"	NR			
5. Acidity	"				
6. PH	"				
7.	"				
8.	"				
9.	"				
10.	"				
11.	"				
12.	"				
13.	"				
14.	"				
15.	"				
16.	"				
17.	"				
18.	"				
19.	"				
20.	"				
21.	"				
22.	"				
23.	"				
Other					
Cyanide					

<sup>1</sup> %R = [ ( SSR - SR ) / SA ] x 100

" N " - out of control

" NR "-Not required

Comments:

## Form M b

Q.C. Report No. 1

## DUPLICATES

LAB NAME Post, Buckley, Schuh & Jernigan  
DATE 4/26/83CASE NO. SAS 3727E  
EPA Sample No. 3727E-06  
Lab Sample ID No. R804014-06  
Units: mg/L

Matrix LA

Compound	Control Limit <sup>1</sup>	Sample (S)	Duplicate (D)	RPD2
1. Sulfate				
2. NO <sub>3</sub> - NO <sub>2</sub>				
3. Chloride				
4. Alkalinity				
5. Acidity		576	578	0.3
6. PH		4.6	4.6	0.0
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
Other				
Cyanide				

\* Out of Control

<sup>1</sup> To be added at a later date.<sup>2</sup> RPD = [ S - D / ((S + D)/2) ] x 100

NC - Non calculable RPD due to value(s) less than CRDL

## Form V C

Q.C. Report No. 1

## DUPLICATES

LAB NAME Post, Buckley, Schuh & Jernigan  
DATE 4/26/88CASE NO. SAS 3727E  
EPA Sample No. 3727E-09  
Lab Sample ID No. 8804014-09  
Units: mg/lMatrix LA

Compound	Control Limit1	Sample (S)	Duplicate (D)	RPD2
1. Sulfate				
2. N03 - N02				
3. Chloride				
4. Alkalinity	+10%	329	32.8	0.3
5. Acidity				
6. PH				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
Other:				
Cyanide				

\* Out of Control

<sup>1</sup> To be added at a later date.<sup>2</sup> RPD = [ (S - D) / ((S + D)/2) ] x 100

NC - Non calculable RPD due to value(s) less than CRDL

Form M a

Q.C. Report No. 1

DUPLICATES

LAB NAME Post, Buckley, Schuh & Jernigan  
DATE 4/26/88

CASE NO. SAS 3727E  
EPA Sample No. 3727 E-01  
Lab Sample ID No. 8804014 - 01  
Units: mg/lc

Matrix LA

Compound	Control Limit <sup>1</sup>	Sample (S)	Duplicate (D)	RPD2
1. Sulfate	<u>±10%</u>	<u>5.88</u>	<u>5.94</u>	<u>1.0</u>
2. NO <sub>3</sub> - NO <sub>2</sub>		<u>0.10</u>	<u>0.10</u>	<u>NC</u>
3. Chloride		<u>0.50</u>	<u>0.50</u>	<u>NC</u>
4. Alkalinity				
5. Acidity				
6. PH				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
Other:				
<u>Cyanide</u>				

\* Out of Control

<sup>1</sup> To be added at a later date.

$$^2 \text{RPD} = [ (S - D) / ((S + D)/2) ] \times 100$$

NC - Non calculable RPD due to value(s) less than CRDL

Form VII  
 Q.C. Report No. 1  
 INSTRUMENT DETECTION LIMITS AND  
 LABORATORY CONTROL SAMPLE  
 LAB NAME Post, Buckley, Schuh, & Jernigan CASE NO SAS 3727E DATE 4/26/66

LCS NO. SAS 3727E

Metals:	Required Detection		Instrument Detection		Lab Control Sample				
	Compound	Limits (CRDL)- <sup>ug/l</sup>	Limits (IDL)- <sup>ug/l</sup>	Furnace	(ug/l)	mg/kg	(circle one)		
		Mg/L		ICP/AA	ID#	ID#	True	Found	%R
Metals:									
1. Sulfate	5.0	5.0			2.0	2.0	100		
2. NO <sub>3</sub> - NO <sub>2</sub>	0.1	0.10			2.0	2.0	100		
3. Chloride	0.5	0.5			52	52	100		
4. Alkalinity	10	10			25	25	100		
5. Acidity	± 10	± 10			6.0	6.0	100		
6. PH	0.1	0.1			6.0	6.0	100		
7.									
8.									
9.									
10.									
11.									
12.									
13.									
14.									
15.									
16.									
17.									
18.									
19.									
20.									
21.									
22.									
23.									
Other									
Cyanide	10		NR	NR					

NR - Not required

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

ESD/Central Regional Laboratory  
DATA TRACKING FORM FOR CONTRACT SAMPLES

CRL Data Set No. SF 5011 CERCLIS No. —  
SMO Case No. SAS 3727E Site Name and Location: Hader Ground Water Contamination  
Name of Contractor or EPA Laboratory: PBS & J Data User: MPCA  
No. of Samples: 9 Date Samples or Data Received: 5-2-88

1. Have chain-of-custody records been received? YES  NO
  2. Have Traffic Reports or packing lists been received? YES  NO
  3. If no, are Traffic Report or packing list numbers written on the chain-of-custody record? YES  NO
  4. If no, which Traffic report or packing list numbers are missing?
- \_\_\_\_\_
- \_\_\_\_\_
- \_\_\_\_\_

Are basic data forms in? YES  NO —

Number of samples claimed: 9 Number of samples received: 9

Checked by: Erinda Lucy M. Aresta Date: 5-2-88

Received by Contract Project Management Section: L.d. Date: 5.3.88

Review Started: 5.11.88 Reviewer Signature: adda heinic

Total time spent on review: 3 Date review completed: 5.12.88

Copied (xeroxed) by: Sylvia Date: 5.12.88

Mailed to Data User by: Pohemes Harris Date: 5-16-88

DATA USERS:

Please fill in the blanks below and return this form to: Sylvia Griffin, Data Management Coordinator, Region V, SSCRL

Data received by: \_\_\_\_\_ Date: \_\_\_\_\_

O.A. review received by: \_\_\_\_\_ Date: \_\_\_\_\_

Inorganic Data Complete [ ]  
Organic Data Complete [ ]  
Dioxin Data Complete [ ]  
SAS Data Complete [ ]  
Suitable for Intended Purposes [ ] ✓ [ ] if acceptable.  
List problems below.

See Attached "Missing Data Request Form" [ ]

PROBLEMS: Please indicate reasons (if any) why data are not suitable for your uses.  
Other problems.

\_\_\_\_\_

\_\_\_\_\_

Received by Data Management Coordinator, CRL for File: Date: \_\_\_\_\_

Signature: \_\_\_\_\_

## Form II.a

Q.C. Report No 1.

## INITIAL AND CONTINUING CALIBRATION VERIFICATIONS

LAB NAME Post, Buckley, Schuh & Jernigan  
 DATE 4/26/81 CASE NO SAS 3727E  
 SOW NO. 7/85  
 UNITS: ~~ppm~~ mg/L

Compound	Initial Calibration1			Continuing Calibration2			Method
	True Value	Found	%R	True Value	Found	%R	
1. Sulfate	8.0	8.6	108	8.0	8.5	106	18.6 108 1375 2
2. N03 - N02	0.8	0.8	100	0.8	0.8	100	0.8 100 1353 2
3. Chloride	4.0	4.0	100	4.3	4.1	102	4.2 105 1335.2
4. Alkalinity	NR						
5. Acidity	1000	992	99	1000	992	99	1305.1
6. PH							
7.							
8.							
9.							
10.							
11.							
12.							
13.							
14.							"
15.							
Other:							
Cyanide							

1 Initial Calib. Source EMSL/LV, 2 Continuing Calib Source Lab prepared3 Control Limits: Mercury and Tin 80-120; Other Metals 90-110;  
Cyanide 85-115

4 Indicate Analytical Method Used: P - ICP; A - Flame; F - Furnace AA

Form II b

Q C Report No 1

INITIAL AND CONTINUING CALIBRATION VERIFICATIONS

LAB NAME Post: Buckley, Schuh & Jernigan  
DATE 4/26/88 CASE NO. SAS 3727E  
SOW NO. 7/85  
UNITS: ~~ug/L~~ mg/L

Compound	Initial Calibration <sup>1</sup>			Continuing Calibration <sup>2</sup>			Method <sup>4</sup>
	True Value	Found	%R	True Value	Found	%R	
1. Sulfate				8.0	8.8	11%	1375.2
2. N03 - N02							
3. Chloride							
4. Alkalinity							
5. Acidity							
6. PH							
7.							
8.							
9.							
10.							
11.							
12.							
13.							
14.							
15.							
Other:							
Cyanide							

1 Initial Calib. Source EMSL/I.V. <sup>No</sup> 2 Continuing Calib. Source Lab prepared

3 Control Limits: Mercury and Tin 80-120; Other Metals 90-110;  
Cyanide 85-115

4 Indicate Analytical Method Used: P - ICP; A - Flame; F - Furnace AA

PBSI

POST, BUCKLEY, SCHUH &amp; JLRNIGAN, INC.

100 NORTH CALIFORNIA  
SUITE 1000  
SAN FRANCISCO, CALIFORNIA 94101  
TELEPHONE 415-362-1100  
FAX 415-362-1111  
TELECO 415-362-1111

88 4/14/88

EPA Method EO 1  
TURBIDITY AND pH DATA SHEET

Corning Model 135

PROJECT NAME: EPA SAs

DATE: 4.2.88

PROJECT NO.: Backup Data

ANALYST: CK

Sample No.	Temperature °C (Temperature probe attached)	pH	Turbidity (NTU)
H.O Buffer		4.001	
T.O Buffer		6.999	
1O.O Buffer		9.954	
WP947	T.V 6.0	5.975	9990
6501 090122	put 3 = 2701- <del>89</del>	6.375	
6501 090119	1044.0-886	6.366	
407 490401401	3727EO1		
501 090120		6.361	Note: Analysis
501 490401401	3727EO1		performed on all sample
501 090120		6.363	
501 490401401	3727EO1		Containers received. (Page 1)
6303 1118 1980101403	3727EO3	7.211	
505 090157		7.269	Since some containers
1123 990101405	3727EO5		are duplicate samples
607 090161		7.046	
132 490401407	3727EO7		one pH is reported
6005 590101409	3727EO9	7.279	on page 2.
142 090149			
104 490401404	3727EO4	6.342	
1434 490401404			
111 090159			
1008 490401408	3727EO8	6.514	
1027 090133			
1503 490401402	3727EO2	6.832	
131 090121			
1501 490401401	3727EO1	6.403	
1501 090121			
1501 490401401	3727EO1	6.391	
1534 090173			
1505 490401406	3727EO6	4.566	
1534 090173			
1505 490401406	3727EO6	4.570	

PBSJ

POST, BUCKLEY, SCHUH & JERNIGAN, INC.

~~100 NORTH DANCE HALL~~  
~~OLANOO ROAD 12001 1004~~  
~~101-111111~~

Page 2

Sept. 9/14/88

EPA method 150.1 Conning pH Meter model 1135

~~TURBIDITY AND PH DATA SHEET~~  
CK 4.2.88

PROJECT NAME: EPA/Sas Case 3727E

DATE: 4.2.88

PROJECT NO.: \_\_\_\_\_

ANALYST: C knepper

Sample No.	Temperature °C Temperature probe attached	pH	Fusibility (HTH) CK 4.2.68
H.C Buffer		4.001	
7.0 Buffer		6.999	
10.0 Buffer		9.954	
WP997	T.V. 6.0	5.975	99%
<u>3727E01</u> <u>990401401</u>		6.383	6.38
<u>3727E02</u> <u>990401402</u>		6.932	6.93
<u>3727E03</u> <u>990401403</u>		7.211	7.21
<u>3727E04</u> <u>990401404</u>		6.342	6.34
<u>3727E05</u> <u>990401405</u>		7.269	7.27
<u>3727E06</u> <u>990401406</u>	$\begin{cases} \% RSD = 0.06 \\ R = 0.004 \end{cases}$	4.566	4.57
<u>3727E06 DUP</u> <u>990401406</u>		4.570	
<u>3727E07</u> <u>990401407</u>		7.048	7.05
<u>3727E08</u> <u>990401408</u>		6.518	6.52
<u>3727E09</u> <u>990401409</u>		7.279	7.28



**POST, DUCKLEY, SCHUM & JERNIGAN, INC.**  
889 NORTH ORANGE AVENUE  
ORLANDO FLORIDA 32801  
305/423 7275

ggy 4/22/88

Corning Potentiometer  
COMPUTER-AIDED TITRIMETER Model 135  
FISHER MODEL 450  
Instrument Log

Case = SAS 3727E

## Buffer

A3:ap



# Post, Buckley, Schuh & Jernigan, Inc.

**CONSULTING ENGINEERS and PLANNERS**

889 NORTH ORANGE AVENUE, ORLANDO, FLORIDA 32801-1088 • 305/423-7276 • TELEX 808435

528 4/22/88

EPA 305.1

## ACIDITY DATA SHEET

PROJECT NAME: SAS 3727E

DATE: 4/6/88

**PROJECT NO.:** \_\_\_\_\_

ANALYST: S.I. Jackson  
F. Boone

Normality of Titrant: 0.0206N HCl or H<sub>2</sub>SO<sub>4</sub> 4/4/88 FB

CALCULATION: Acidity as mg/l CaCO<sub>3</sub> =  $\frac{(A \times B) - (C \times D) \times 5 \times 10^4}{\text{ml sample}}$

Normality NaOH =

$$\frac{A \times B}{204.2 \times C}$$

A = Volume of std. NaOH used intitration

B = Normality of std. NaOH = 0.0190N 4/6/88 FB

C = Volume of std H<sub>2</sub>SO<sub>4</sub> used to reduce pH to 4 or less

D = Normality of std.  $\text{H}_2\text{SO}_4$

$$A = g \cdot KNP / l = 10.09$$

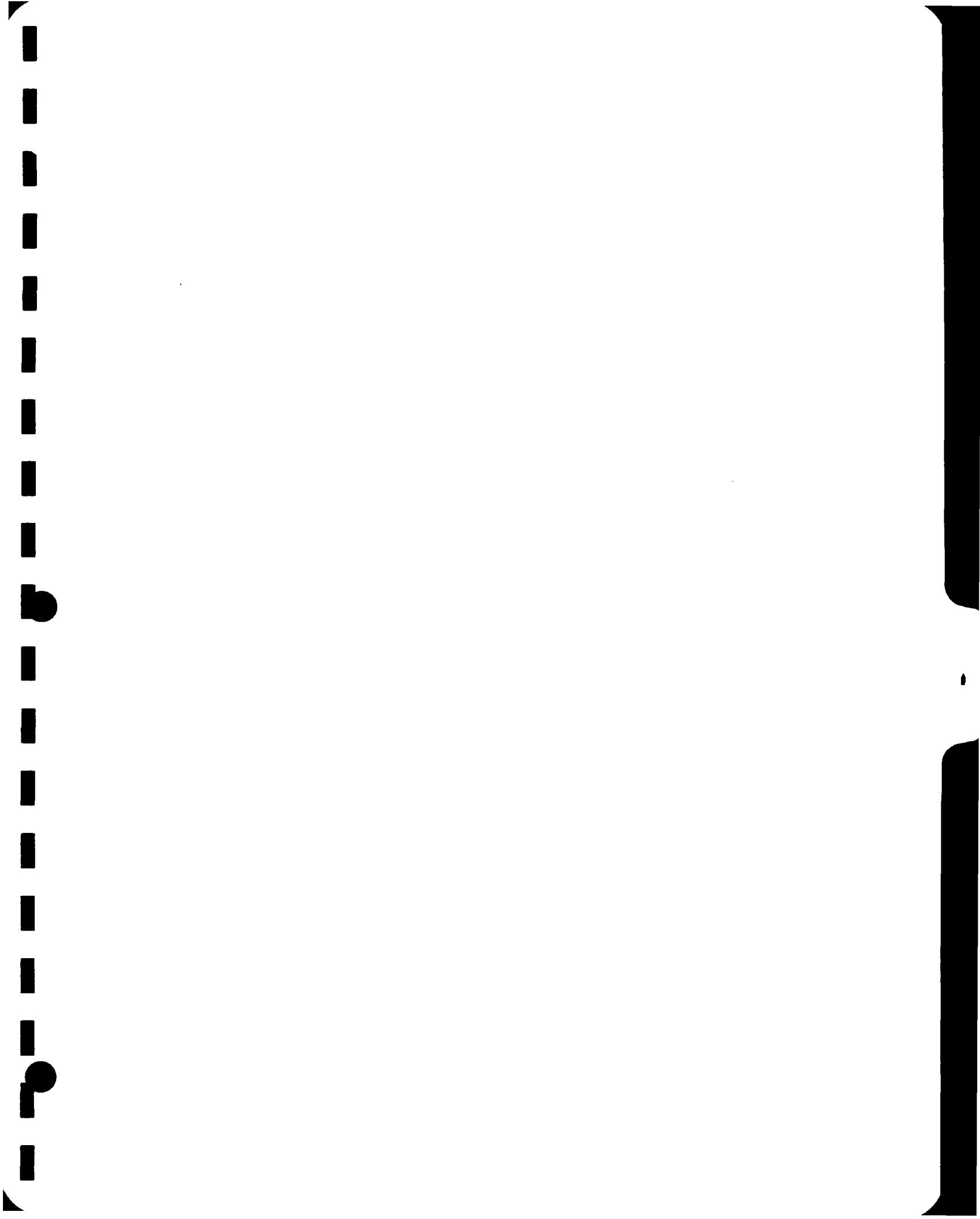
$$\text{Normality } H_2SO_4 = \frac{A \times B}{53.0 \times C}$$

B- abs KHP titrated = 10.0 abs

$$A = g \text{Nb}_2\text{O}_3/\text{l} = 2.59/\text{l}$$

$$g = \text{mols NaOH used} = 25.9 \text{ mols}$$

$$C = \text{molar } H_2SO_4 \text{ used} = 4.57 \text{ M}$$



Dissolved!

MINNESOTA DEPARTMENT OF HEALTH  
ANALYTICAL DATA

BILLING NO:

MPCA

032

Samples Collected By Ruotsinoja / Adams

Report To Shawn Ruotsinoja

Field Number	River, Town, Etc.	Sampling Point and Source of Sample
2 a	Hader	non responsive
4 b	Hader	
6 c	Hader	
8 d	Hader	
10 e	Hader	

This line for Lab use only

Sample Number

132241

a 132242

b 132243

c 132244

d 132245

Date Collected 8/5/87

Time Collected

Date Received by Lab 8-17-87

Temperature, °C

Dissolved Oxygen

pH value, SU 013

Total Residual Chlorine

Fecal Coliform, MPN/100 ml 305

Fecal Strep., No/100 ml 313

Total Solids 001

Suspended Solids 003

Turbidity, NTU 011

Calcium as CaCO<sub>3</sub> 251 (252)

210.

250.

380.

220.

190.

Magnesium as CaCO<sub>3</sub> 253 (254)

100.

110.

170.

110.

90.

Total Hardness as CaCO<sub>3</sub> 021

Chloride as Cl 023

5-Day B.O.D. 096

Nitrification Inhibited BOD<sub>5</sub> 083

Total Phosphorus as P 059

Orthophosphorus as P 063

Organic Nitrogen as N 065

Ammonia Nitrogen as N 064

Nitrite + Nitrate Nitro. 069

Nitrite Nitrogen as N 067

Arsenic as As 108

Cadmium as Cd \*125 (123)

0.035

0.038

0.012

0.010

0.017

Total Chromium as Cr 129

Hexavalent Chromium as Cr 034

Copper as Cu 145/143

Iron as Fe 152 (151)

6700.

22.

21000.

2500.

6200.

Lead as Pb 159 (158)

<0.2

0.7

<0.2

<0.2

<0.2

Manganese as Mn 166 (165)

59.

13.

440.

74.

26.

Mercury as Hg 200

Nickel as Ni 173 (172)

16.

<1.0

9.9

<1.0

9.2

Zinc as Zn 194 (193)

500.

25.

60.

120.

440.

Flame Metals Ashing Code 207

Phenol µg/l 085

Oil and Grease 089

Chlorophyll a, µg/l 450

Kieldehl Nitrogen 068

Flame/Furnace

RECEIVED

COMPLETED

DEC 9 1987

MPCA, HAZARDOUS  
WASTE DIVISION

CHEMICAL LABS

MINNESOTA DEPARTMENT OF HEALTH  
SUPPLEMENTAL SHEET

## **ANALYTICAL DATA**

Samples Collected By Ruotsinaja / Adams

Report To Shawn Ruotsaloja

**RECEIVED**

~~DEC 9 1987~~

~~MPCA, HAZARDOUS  
WASTE DIVISION~~

**COMPLETED**

DEC 07 1987

CHEMICAL LABS

- 1 -

Dissolved

MINNESOTA DEPARTMENT OF HEALTH  
ANALYTICAL DATA

BILLING NO:

MPCA

032

Samples Collected By Ruotsinoja / Adams

Report To ✓ Shawn Ruotsinoja

Field Number	River, Town, Etc.	Sampling Point and Source of Sample
12 a	Hader	non responsive
14 b	"	
16 c	"	
18 d	"	
20 e	"	

This line for Lab use only

Sample Number	132246 a	132247 b	132249 c	132249 d	132250 e
---------------	----------	----------	----------	----------	----------

Date Collected 8/6/87

Time Collected

Date Received by Lab 8-17-87

Temperature, °C

Dissolved Oxygen

pH value, SU 013

Total Residual Chlorine

Fecal Coliform, MPN/100 ml 305

Fecal Strep., No/100 ml 313

Total Solids 001

Suspended Solids 003

Turbidity, NTU 011

Calcium as CaCO<sub>3</sub>

252 220. 430. 300. 260. 190.

Magnesium as CaCO<sub>3</sub>

253 254) 110. 170. 140. 140. 80.

Total Hardness as CaCO<sub>3</sub> 021

Chloride as Cl

5 Day B.O.D. 096

Nitrification Inhibited BOD<sub>5</sub> 083

Total Phosphorus as P 059

Orthophosphorus as P 063

Organic Nitrogen as N 065

Ammonia Nitrogen as N 064

Nitrite + Nitrate Nitro. 069

Nitrite Nitrogen as N 067

Arsenic as As 108

Cadmium as Cd \*125/129 123 <0.010

Total Chromium as Cr 129

Hexavalent Chromium as Cr 034

Copper as Cu 145/143

Iron as Fe 152/150 151 2900. 210000. 93000. 690. 80000.

Lead as Pb 159/157 158 0.4 45. 1.1 20.2 2.8

Manganese as Mn 166/164 164 44. 3000. 500. 130. 600.

Mercury as Hg 200

Nickel as Ni 173/172 172 <1. 390. 220. 3.0 170.

Zinc as Zn 194/193 193 490. 3200. 1000. 10. 1330.

Flame Metals Ashing Code 207

Phenol µg/l 085

Oil and Grease 089

Chlorophyll a, µg/l 450

Kjeldahl Nitrogen 068

Flame/Furnace

RECEIVED

COMPLETED

DEC 07 1987

DEC 9 1987

CHEMICAL LABS

MPCA, HAZARDOUS  
WASTE DIVISION

MINNESOTA POLLUTION CONTROL AGENCY  
SUPPLEMENTAL SHEET

## **ANALYTICAL DATA**

MPCA

Samples Collected By Hufschmid / Adams

Report To Shawn Ruotsinoja

Results are in milligrams per liter except as otherwise noted.

## MINNESOTA DEPARTMENT OF HEALTH

## ANALYTICAL DATA

BILLING NO:

- 395

Samples Collected By Ruotsingja / AdamsReport To Shawn Ruotsingja  
Byron Adams

Field Number	River, Town, Etc.	Sampling Point and Source of Sample
A a	Hader	non responsive
B b	"	
C c	"	
D d	"	
E e	"	

This Line for Lab use only				
Sample Number	132251	a 132252	b 132253	c 132254
Date Collected	8/5/87			d 132255
Time Collected				e
Date Received by Lab	8-17-87			
Temperature, °C				
Dissolved Oxygen				
pH Value, SU	013			
Total Residual Chlorine				
total Coliform, MPN/100 ml	305			SEP 26 1987
total Strep., No/100 ml	313			JFH MPCA SOLID & HAZ WASTE DIVISION
Total Solids	001			
Suspended Solids	003			
Transparency, NTU	011			
Calcium as CaCO <sub>3</sub>				
Magnesium as CaCO <sub>3</sub>				
Total Hardness as CaCO <sub>3</sub>	021			
Chloride as Cl	023	<0.50	29.	0.38
-Day B.O.D.	096			<0.50
Identification Inhibited BOD <sub>5</sub>	083			<0.50
Total Phosphorus as P	059			
Inorganic Phosphorus as P	063			
Organic Nitrogen as N	065			
Inorganic Nitrogen as N	064			
Chloride + Nitrate Nitro.	069			
Nitrite Nitrogen as N	067			
Arsenic as As	108			
Cadmium as Cd	*125/122			
Total Chromium as Cr	129			
Hexavalent Chromium as Cr	034			
Copper as Cu	145/143			
Iron as Fe	152/150			
Lead as Pb	159/157			
Manganese as Mn	166/164			
Mercury as Hg	200			
Nickel as Ni	173/171			
Zinc as Zn	194/192			
Trace Metals Ashing Code	207			
Iron µg/l	085			
Oil Grease	089			
Chlorophyll a, µg/l	450			
Lehman Nitrogen	068			
Lake/Furnace				

COMPLETED

SEP 21 1987

CHEMICAL LABS

**MINNESOTA DEPARTMENT OF HEALTH  
SUPPLEMENTAL SHEET**

## **ANALYTICAL DATA**

Samples Collected By Ruotsinaja / Adams

Report To Shawn Ruotsinga  
Byron Adams

Results are in milligrams per liter except as otherwise noted.

## MINNESOTA POLLUTION CONTROL AGENCY

## ANALYTICAL DATA

BILLING NO: PCA - 032

MPCA

Samples Collected By Ruotsincja / AdamsReport To Shawn Ruotsincja  
Byron Adams

Field Number	River, Town, Etc.	Sampling Point and Source of Sample
F a	Hader	non responsive
G b	"	
I c	"	
K d	"	
L e	"	

This line for Lab use only					
Sample Number	132256 a	132257 b	132258 c	132259 d	132260 e
Date Collected	8/6/87				
Time Collected					
Date Received by Lab	8-17-87				
Temperature, °C					
Dissolved Oxygen					
pH value, SU	013				
Total Residual Chlorine					
Fecal Coliform, MPN/100 ml	305				
Fecal Strep., No/100 ml	313				
Total Solids	001				
Suspended Solids	003				
Turbidity, NTU	011				
Calcium as CaCO <sub>3</sub>	251				
Magnesium as CaCO <sub>3</sub>	253				
Total Hardness as CaCO <sub>3</sub>	021				
Chloride as Cl	(023)	<0.50	<0.50	<0.50	7.7
5-Day B.O.D.	096				
Nitrification Inhibited BOD <sub>5</sub>	083				
Total Phosphorus as P	059				
Orthophosphorus as P	063				
Organic Nitrogen as N	065				
Ammonia Nitrogen as N	064				
Nitrite + Nitrate Nitro.	069				
Nitrite Nitrogen as N	067				
Arsenic as As	108				
Cadmium as Cd	*125/122				
Total Chromium as Cr	129				
Hexavalent Chromium as Cr	034				
Copper as Cu	145/143				
Iron as Fe	152/150				
Lead as Pb	159/157				
Manganese as Mn	166/164				
Mercury as Hg	200				
Nickel as Ni	173/171				
Zinc as Zn	194/192				
Flame Metals Ashing Code	207	JULY 1987	2017		
Phenol µg/l	085	JULY 1987	2017		
Oil and Grease	089	JULY 1987	2017		
Chlorophyll a, µg/l	450	JULY 1987	2017		
Tjeldahl Nitrogen	068	JULY 1987	2017		

#Flame/Furnace

COMPLETED

SEP 21 1987

CHEMICAL LABS

**M**INNESOTA POLLUTION CONTROL AGENCY  
SUPPLEMENTAL SHEET

## **ANALYTICAL DATA**

Samples Collected By Ruktineja / Adams

Report To Shawn Ruotsalo  
Byron Adams

Results are in milligrams per liter except as otherwise noted.

## MINNESOTA DEPARTMENT OF HEALTH

## ANALYTICAL DATA

BILLING NO:

MPGA

032

Samples Collected By Ruotsinaja/AdamsReport To Shawn Ruotsinaja

Field Number	River, Town, Etc.	Sampling Point and Source of Sample
1 a	Hader	non responsive
3 b	Hader	
5 c	Hader	
7 d	Hader	
9 e	Hader	

This line for Lab use only	a	b	c	d	e	
Sample Number	132231	132232	132233	132234	132235	
Date Collected	8/5/87					
Time Collected						
Date Received by Lab	8-17-87					
Temperature, °C						
Dissolved Oxygen						
pH value, SU	013					
Total Residual Chlorine						
Fecal Coliform, MPN/100 ml	305					
Fecal Strep., No/100 ml	313					
Total Solids	001					
Suspended Solids	003					
Turbidity, NTU	011					
Calcium as CaCO <sub>3</sub>	251					
Magnesium as CaCO <sub>3</sub>	253					
Total Hardness as CaCO <sub>3</sub>	021					
Chloride as Cl	023					
5 Day B.O.D.	096					
Nitrification Inhibited BOD <sub>5</sub>	083					
Total Phosphorus as P	059					
Orthophosphorus as P	063					
Organic Nitrogen as N	065					
Ammonia Nitrogen as N	064					
Nitrite + Nitrate Nitro.	069					
Nitrite Nitrogen as N	067					
Arsenic as As	108					
Cadmium as Cd	*125/122	0.016	0.014	0.010	0.057	0.020
Total Chromium as Cr	129					
Hexavalent Chromium as Cr	034					
Copper as Cu	145/143					
Iron as Fe	152/150	690.	23.	100.	6400.	380.
Lead as Pb	159/157	0.8	0.9	0.6	1.3	0.5
Manganese as Mn	166/164	28.	9.	15.	82.	<3.
Mercury as Hg	200					
Nickel as Ni	173/171	17.	<1.0	<1.0	<1.0	2.3
Zinc as Zn	194/192	15.	11.	6.1	1600.	9.8
Metals Ashing Code	207					
Total µg/l	085					
and Grease	089					
Prophylia, µg/l	450					
Dissolved Nitrogen	068					
Amine/Furnace						

COMPLETED

SEP 04 1987

CHEMICAL LABS

RECEIVED

SEP 9 1987

MPGA, SOLID & HAZ  
WASTE DIVISION

5/11

## MINNESOTA DEPARTMENT OF HEALTH

## ANALYTICAL DATA

BILLING NO:

032

Samples Collected By Ruotsinaja / Adams

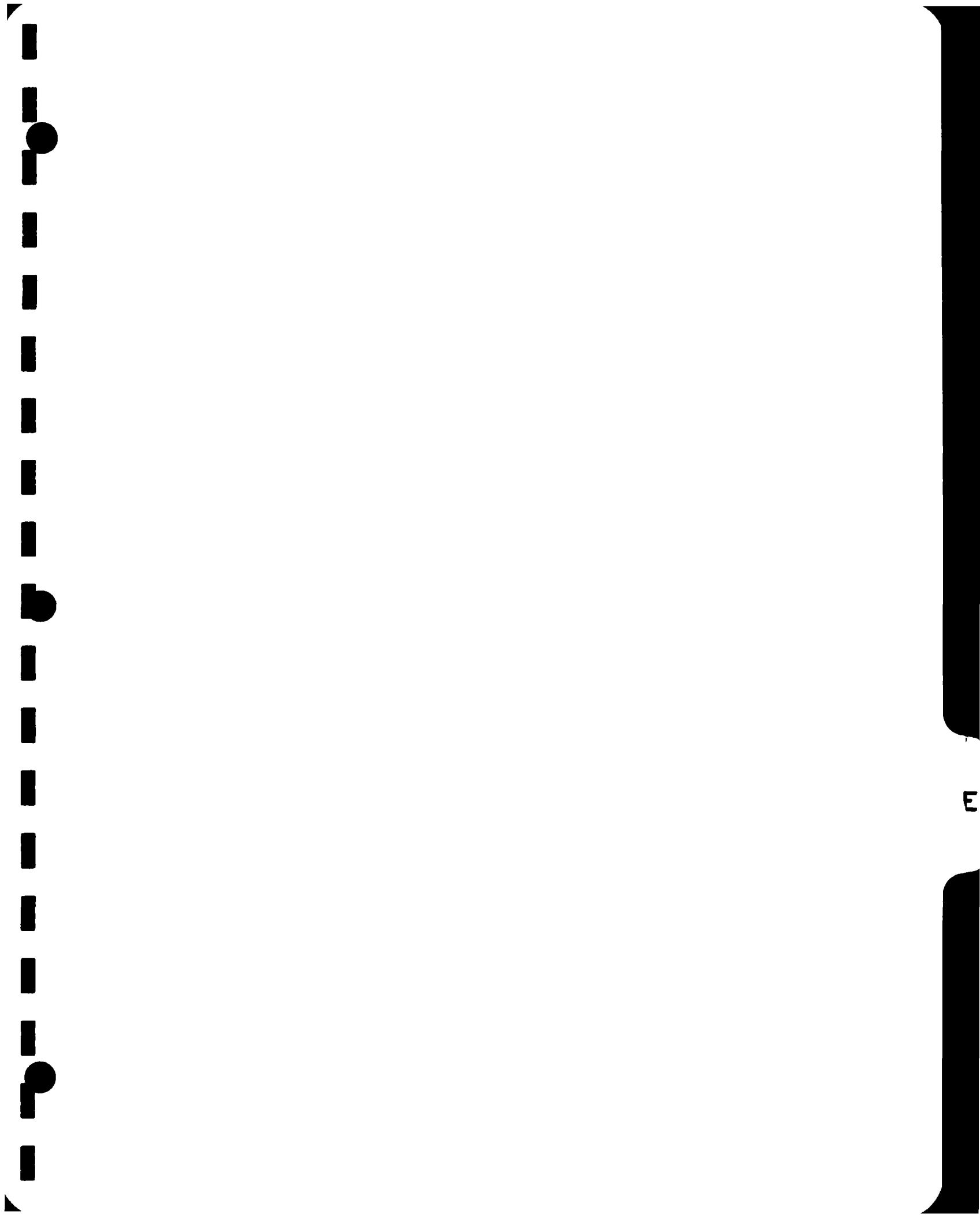
Report To

Shawn Ruotsinaja

MPCA

Field Number	River, Town, Etc.	Sampling Point and Source of Sample				
11 a	Hader		non responsive			
13 b	Hader					
15 c	Hader					
17 d	Hader					
19 e	Hader					

This line for Lab use only	a	b	c	d	e	
Sample Number	132236	132237	132238	132239	132241	
Date Collected	8/6/87					
Time Collected						
Date Received by Lab	8-17-87					
Temperature, °C						
Dissolved Oxygen						
pH value, SU	013					
Total Residual Chlorine						
Fecal Coliform, MPN/100 ml	305					
Fecal Strep., No/100 ml	313					
Total Solids	001					
Suspended Solids	003					
Turbidity, NTU	011					
Calcium as CaCO <sub>3</sub>	251					
Magnesium as CaCO <sub>3</sub>	253					
Total Hardness as CaCO <sub>3</sub>	021					
Chloride as Cl	023					
Day B.O.D.	096					
Nitrification Inhibited BOD <sub>5</sub>	083					
Total Phosphorus as P	059					
Orthophosphorus as P	063					
Organic Nitrogen as N	065					
Ammonia Nitrogen as N	064					
Nitrite + Nitrate Nitro.	069					
Nitrite Nitrogen as N	067					
Arsenic as As	108					
Cadmium as Cd	*125/122	<0.010	0.010	0.42	0.019	0.010
Total Chromium as Cr	129					
Hexavalent Chromium as Cr	034					
Copper as Cu	145/143					
Iron as Fe	152/150	240.	51.	140000.	1400.	930.
Lead as Pb	159/157	0.2	0.9	22.	0.2	46.
Manganese as Mn	166/164	3.	3.	650.	160.	9.
Mercury as Hg	200					
Nickel as Ni	173/171	<1.0	<1.0	350.	2.7	24.
Zinc as Zn	194/192	18.	24.	1500.	260.	36.
Flame Metals Ashing Code	207					
Phenol µg/l	085					
Oil and Grease	089					
Chlorophyll a, µg/l	450					
Weldahl Nitrogen	068					
#Flame/Furnace						



**STORED RETRIEVAL DATE 87/07/29**

PGM=INVENT

442128092570701 1101802CDCA1 14468  
44 21 28.0 .092 .57 07.0 2  
AQF: PRAIRIE DU CHIEN /  
27049 MINNESOTA GOODHUE  
MAJOR BASIN: UPPER MISS 070639  
MINOR BASIN: LOWER UPPER MISS RIVER  
21MINNG 821211 070400  
0000 FEET DEPTH

PAGE:

GW0035

367PRDX

/TYPE/AMOUNT/WELL

STORET RETRIEVAL DATE 87/07/29

PGM=INVENT

PAGE: 2  
GWQ0351

442128092570701 1101802CDCA1 144680  
 44 21 28.0 092-57 07.0 /2  
 AQF: PRAIRIE DU CHIEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070639  
 MINOR BASIN: LOWER UPPER MISS RIVER  
 21MINNG 821211 07040002 HQ  
 0000 FEET DEPTH

/TYP/A/MBNT/WELL

PARAMETER		MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
01092 ZINC	ZN, TOT	UG/L	WATER		1 70.00000			70	70	82/06/29	82/06/29
01147 SELENIUM	SE, TOT	UG/L	WATER	K	1 1.000000			1	1	82/06/29	82/06/29
31505 TOT COLI	MPN CONF	/100ML	WATER	K	1 2.000000			2	2	82/10/14	82/10/14
				L	1 24000.00			24000	24000	82/06/29	82/06/29
			TOT	2	12001.00	2879E+05	16969.00	24000	2	82/06/29	82/10/14
31615 FEC COLI	MPNECMED	/100ML	WATER	K	1 2.000000			2	2	82/10/14	82/10/14
31679 FECSTREP MF M-ENT		/100ML	WATER	K	1 2.000000			2	2	82/10/14	82/10/14
32101 DICLBRMT		TOTUG/L	WATER	K	1 .5000000			.5	.5	86/10/02	86/10/02
32102 CARBNET		TOTUG/L	WATER	K	1 .2000000			.2	.2	86/10/02	86/10/02
32104 BROMOFRM	WHL-WTR	UG/L	WATER	K	1 1.000000			1.0	1.0	86/10/02	86/10/02
32106 CHLRFORM		TOTUG/L	WATER	K	1 .2000000			.2	.2	86/10/02	86/10/02
32730 PHENOLS	TOTAL	UG/L	WATER	K	1 2.000000			2	2	82/06/29	82/06/29
34010 TOLUENE		TOT UG/L	WATER	K	1 .5000000			.50	.50	86/10/02	86/10/02
34030 BENZENE		TOT UG/L	WATER	K	1 .5000000			.50	.50	86/10/02	86/10/02
34301 CHLOROBE	NZENE	TOTWUG/L	WATER	K	1 .5000000			.500	.500	86/10/02	86/10/02
34306 CHLORODI	BROMOMET	TOTWUG/L	WATER	K	1 .5000000			.500	.500	86/10/02	86/10/02
34371 ETHYLBEN	ZENE	TOTWUG/L	WATER	K	1 .5000000			.500	.500	86/10/02	86/10/02
34423 METHYLEN	ECHLORID	TOTWUG/L	WATER	K	1 1.000000			1.000	1.000	86/10/02	86/10/02
34475 TETRACHL	OROETHYL	TOTWUG/L	WATER	K	1 2.000000			2.000	2.000	86/10/02	86/10/02
34488 TRICHLOR	OFLUOROM	TOTWUG/L	WATER	K	1 .5000000			.500	.500	86/10/02	86/10/02
34496 11DICHLO	ROETHANE	TOTWUG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
34501 11DICHLO	ROETHYLE	TOTWUG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
34506 111TRICH	LOROETHA	TOTWUG/L	WATER	K	1 .5000000			.500	.500	86/10/02	86/10/02
34511 112TRICH	LOROETHA	TOTWUG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
34516 1122TETR	ACHLOROE	TOTWUG/L	WATER	K	1 2.000000			2.000	2.000	86/10/02	86/10/02
34531 12DICHLO	ROETHANE	TOTWUG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
34536 12DICHLO	ROBENZEN	TOTWUG/L	WATER	K	1 1.000000			1.000	1.000	86/10/02	86/10/02
34541 12DICHLO	ROPROPAN	TOTWUG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
34546 12DICHLO	ROETHENE	TOTWUG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
34566 13DICHLO	ROBENZEN	TOTWUG/L	WATER	K	1 1.000000			1.000	1.000	86/10/02	86/10/02
34571 14DICHLO	ROBENZEN	TOTWUG/L	WATER	K	1 1.000000			1.000	1.000	86/10/02	86/10/02
34699 T1,3-DCP	TOT WAT	UG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
34704 C1,3-DCP	TOT WAT	UG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
39180 TRICHLOR	ETHYLENE	TOT UG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
70300 RESIDUE	DISS-180	C MG/L	WATER	K	1 310.0000			310	310	82/06/29	82/06/29
71900 MERCURY	HG, TOTAL	UG/L	WATER	K	1 .1000000			.1	.1	82/06/29	82/06/29
74041 WQF	SAMPLE	UPDATED	WATER	2	865670.0	41943000	6476.400	870206	861125	86/10/02	86/10/02
77093 C-1,2DCE	TOTAL	UG/L	WATER	K	1 .2000000			.200	.200	86/10/02	86/10/02
77134 M-XYLENE	TOTAL	UG/L	WATER	K	1 .5000000			.500	.500	86/10/02	86/10/02

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PGM=INVENT

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GWQ0351

442128092570701 1101802CDCA1 144680

44. 21. 28. 0. 092-57. 07. 0 2

AQF: PRAIRIE DU CHIEN /

367PRDC

27049 MINNESOTA GOODHUE

MAJOR BASIN: UPPER MISS 070639

MINOR BASIN: LOWER UPPER MISS RIVER

21MINNG 821211 07040002 HQ

0000 FEET DEPTH

/TYP/A/MBNT/WELL

	PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
77166	2,3DCLPR TOTAL	UG/L WATER	K	1	.2000000			.200	.200	86/10/02	86/10/02
77168	1,1DCLPR TOTAL	UG/L WATER	K	1	.2000000			.200	.200	86/10/02	86/10/02
77223	IPROPBNZ TOTAL	UG/L WATER	K	1	1.000000			1.000	1.000	86/10/02	86/10/02
77596	DBRMETHA TOTAL	UG/L WATER	K	1	1.000000			1.000	1.000	86/10/02	86/10/02
77651	1,2DBRET TOTAL	UG/L WATER	K	1	.5000000			.500	.500	86/10/02	86/10/02
77652	112C122F TOTAL	UG/L WATER	K	1	.2000000			.200	.200	86/10/02	86/10/02
78109	ALLYLCLR TOT WH W	UG/L WATER	K	1	.5000000			.50	.50	86/10/02	86/10/02
78121	P-XYLEN+ O-XYLEN	TOT UG/L WATER	K	1	.5000000			.50	.50	86/10/02	86/10/02
81552	ACETONE	TOT UG/L WATER	K	1	20.00000			20.000	20.000	86/10/02	86/10/02
81578	DIETHYL ETHER	TOT UG/L WATER	K	1	1.000000			1.000	1.000	86/10/02	86/10/02
81595	MTH ETH KETONE	TOT UG/L WATER	K	1	5.000000			5.000	5.000	86/10/02	86/10/02
81596	MTHISOBU KETONE	TOT UG/L WATER	K	1	2.000000			2.000	2.000	86/10/02	86/10/02
81607	TETRAHYD FURAN	TOT UG/L WATER	K	1	10.00000			10.000	10.000	86/10/02	86/10/02
81611	TRICL TRIFLETH	TOT UG/L WATER	K	1	.5000000			.500	.500	86/10/02	86/10/02
84000	GEOLOGIC AGE	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84001	AQUIFER NAME	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84002	CODE GENERAL	REMARKS WATER	TXT	4	TEXT	TEXT	TEXT	TEXT	TEXT	82/06/29	86/10/02

STORET RETRIEVAL DATE 87/07/29

PGM=INVENT

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GWQ0137

442003092423401 1101614BBDD1 144668  
 44 20 12.1 092 42 21.6 2  
 AQF: PRAIRIE DU CHIEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070641  
 MINOR BASIN: LOWER MISS RIVER  
 21MINNG 790122 07040004  
 0000 FEET DEPTH

/TYP/A/AMBN/T/WELL

PARAMETER	LAB IDENT.	NUMBER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
00005	LAB	IDENT.	WATER		2	125720.0	.0000000	.0000000	125716	125715	83/10/06	83/10/06
00010	WATER	TEMP	CENT		2	9.500000	.0000000	.0000000	9.5	9.5	78/10/25	83/10/06
00080	COLOR	PT-CO	UNITS		1	5.000000			5	5	78/10/25	78/10/25
00094	CNDUCTVY	FIELD	MICROMHO		2	478.5000	420.5000	20.50600	493	464	78/10/25	83/10/06
00095	CNDUCTVY	AT 25C	MICROMHO		1	500.0000			500	500	78/10/25	78/10/25
00136	SAMPLE TEMP AT	LAB DEGC	WATER		1	2.000000			2.00000	2.00000	83/10/06	83/10/06
00400	PH	SU	WATER		2	7.150000	.2450000	.4949700	7.50	6.80	78/10/25	83/10/06
00403	LAB	PH	SU		2	7.500000	.0000000	.0000000	7.5	7.5	78/10/25	83/10/06
00410	T ALK	CACO3	MG/L		2	255.0000	50.00000	7.071100	260	250	78/10/25	83/10/06
00425	HCO3 ALK	CACO3	MG/L		2	255.0000	50.00000	7.071100	260	250	78/10/25	83/10/06
00431	T ALK	FIELD	MG/L		2	261.0000	162.0000	12.72800	270	252	78/10/25	83/10/06
00505	RESIDUE TOT VOL	MG/L	WATER	K	1	47.00000			47	47	78/10/25	78/10/25
00615	NO2-N TOTAL	MG/L	WATER	K	1	.0100000			.010	.010	78/10/25	78/10/25
00625	TOT KJEL N	MG/L	WATER		1	.1700000			.170	.170	78/10/25	78/10/25
00630	NO2&NO3 N-TOTAL	MG/L	WATER		2	.0250000	.0004500	.0212130	.04	.01	78/10/25	83/10/06
00665	PHOS-TOT	MG/L P	WATER	K	1	.0010000			.001	.001	78/10/25	78/10/25
00680	T ORG C C	MG/L	WATER	K	2	1.000000	.0000000	.0000000	1.0	1.0	78/10/25	83/10/06
00900	TOT HARD	CACO3	MG/L		2	265.0000	50.00000	7.071100	270	260	78/10/25	83/10/06
00910	CALCIUM	CACO3	MG/L		2	175.0000	50.00000	7.071100	180.0	170.0	78/10/25	83/10/06
00920	MGSNMIUM	CACO3	MG/L		2	92.00000	8.000000	2.828400	94.0	90.0	78/10/25	83/10/06
00929	SODIUM NA,TOT	MG/L	WATER		1	2.300000			2.30	2.30	83/10/06	83/10/06
00930	SODIUM NA,DISS	MG/L	WATER		1	2.100000			2.10	2.10	78/10/25	78/10/25
00935	PTSSIUM K,DISS	MG/L	WATER		1	1.000000			1.00	1.00	78/10/25	78/10/25
00937	PTSSIUM K,TOT	MG/L	WATER		1	1.300000			1.30	1.30	83/10/06	83/10/06
00940	CHLORIDE TOTAL	MG/L	WATER	K	1	.6800000			.7	.7	83/10/06	83/10/06
				K	1	.5000000			.5	.5	78/10/25	78/10/25
				TOT	2	.5900000	.0162010	.1272800	.7	.5	78/10/25	83/10/06
00945	SULFATE SO4-TOT	MG/L	WATER		1	16.00000			16	16	83/10/06	83/10/06
00946	SULFATE SO4-DISS	MG/L	WATER		1	17.00000			17.0	17.0	78/10/25	78/10/25
00950	FLUORIDE F,DISS	MG/L	WATER		1	.1400000			.14	.14	78/10/25	78/10/25
00955	SILICA DISOLVED	MG/L	WATER		1	13.00000			13.0	13.0	78/10/25	78/10/25
01020	BORON B,DISS	UG/L	WATER	K	1	50.00000			50	50	78/10/25	78/10/25
01027	CADMUM CD,TOT	UG/L	WATER		1	.0260000			.03	.03	78/10/25	78/10/25
01034	CHROMIUM CR,TOT	UG/L	WATER	K	1	.5000000			.5	.5	78/10/25	78/10/25
01045	IRON FE,TOT	UG/L	WATER		2	260.0000	3200.000	56.56900	300	220	78/10/25	83/10/06
01051	LEAD PB,TOT	UG/L	WATER		1	2.400000			2	2	78/10/25	78/10/25
01055	MANGNESE MN	UG/L	WATER		2	40.00000	200.0000	14.14200	50.0	30.0	78/10/25	83/10/06
01067	NICKEL NI,TOTAL	UG/L	WATER		1	4.600000			5	5	78/10/25	78/10/25
01092	ZINC ZN,TOT	UG/L	WATER		1	240.00000			240	240	78/10/25	78/10/25

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PGM-INVENT

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/TYP/A/MBNT/WELL

442003092423401 1101614BBDD1 144668  
 44 20 12.1 092 42-21.6 2  
 AQF: PRAIRIE DU CHIEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070641  
 MINOR BASIN: LOWER MISS RIVER  
 21MINNG 790122 07040004  
 0000 FEET DEPTH

GWQ0137

367PRDC

PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
31505 TOT COLI MPN CONF	/100ML WATER		1	2.000000			2	2	78/10/25	78/10/25
		K	1	2.200000			2	2	83/10/06	83/10/06
		TOT	2	2.100000	.0200030	.1414300	2	2	78/10/25	83/10/06
31615 FEC COLI MPNECMED	/100ML WATER	K	1	2.000000			2	2	78/10/25	78/10/25
32101 DICLBRMT	TOTUG/L WATER	K	1	.5000000			.5	.5	83/10/06	83/10/06
32102 CARBTET	TOTUG/L WATER	K	1	.2000000			.2	.2	83/10/06	83/10/06
32104 BROMOFRM WHL-WTR	UG/L WATER	K	1	1.000000			1.0	1.0	83/10/06	83/10/06
32106 CHLRFORM	TOTUG/L WATER	K	1	.2000000			.2	.2	83/10/06	83/10/06
34010 TOLUENE	TOT UG/L WATER	K	1	.5000000			.50	.50	83/10/06	83/10/06
34030 BENZENE	TOT UG/L WATER	K	1	.5000000			.50	.50	83/10/06	83/10/06
34301 CHLOROBENZENE	TOTWUG/L WATER	K	1	.5000000			.500	.500	83/10/06	83/10/06
34306 CHLORODIBROMOMET	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/06	83/10/06
34371 ETHYLBENZENE	TOTWUG/L WATER	K	1	.5000000			.500	.500	83/10/06	83/10/06
34423 METHYLENECHLORID	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/06	83/10/06
34475 TETRACHLOROETHYL	TOTWUG/L WATER	K	1	2.000000			2.000	2.000	83/10/06	83/10/06
34480 THALLIUM SEDMG/KG	DRY WGT WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34496 11DICHLO ROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34501 11DICHLO ROETHYLE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34506 111TRICHLOROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34511 112TRICHLOROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34516 112TETRACHLOROETHANE	TOTWUG/L WATER	K	1	2.000000			2.000	2.000	83/10/06	83/10/06
34531 12DICHLO ROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34536 12DICHLO ROBENZEN	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/06	83/10/06
34541 12DICHLO ROPROPA	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34546 12DICHLO ROETHENE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34566 13DICHLO ROBENZEN	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/06	83/10/06
34571 14DICHLO ROBENZEN	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/06	83/10/06
34576 2CHLOROETHYL VINYL	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/06	83/10/06
34699 T1,3-DCP TOT WAT	UG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
34704 C1,3-DCP TOT WAT	UG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
39180 TRICHLOROETHYLENE	TOT UG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
70300 RESIDUE DISS-180 C	MG/L WATER		1	270.0000			270	270	78/10/25	78/10/25
71900 MERCURY HG,TOTAL	UG/L WATER		1	.4200000			.4	.4	78/10/25	78/10/25
77093 C-1,2DCE TOTAL	UG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
77134 M-XYLENE TOTAL	UG/L WATER	K	1	.5000000			.500	.500	83/10/06	83/10/06
77135 O-XYLENE TOTAL	UG/L WATER	K	1	.5000000			.500	.500	83/10/06	83/10/06
77166 2,3DCLPR TOTAL	UG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
77168 1,1DCLPR TOTAL	UG/L WATER	K	1	.2000000			.200	.200	83/10/06	83/10/06
77173 1,3DCLPR TOTAL	UG/L WATER	K	1	3.000000			3.000	3.000	83/10/06	83/10/06

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442003092423401 1101614BBDD1 144668  
 44 20 12.1 092 42 21-6 2  
 AQF: PRAIRIE DU CHIEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070641  
 MINOR BASIN: LOWER UPPER MISS RIVER  
 21MINNG 790122 07040004  
 0000 FEET DEPTH

GWQ0137

367PRDC

/TYP/A/MBNT/WELL

	PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
77223	IPROPBNZ TOTAL	UG/L WATER	K	1	.5000000		.500	.500	83/10/06	83/10/06	
77443	1,2,3TCP TOTAL	UG/L WATER	K	1	2.000000		2.000	2.000	83/10/06	83/10/06	
77562	1112TCLE TOTAL	UG/L WATER	K	1	.2000000		.200	.200	83/10/06	83/10/06	
77596	DBRMETHA TOTAL	UG/L WATER	K	1	1.000000		1.000	1.000	83/10/06	83/10/06	
77651	1,2DBRET TOTAL	UG/L WATER	K	1	1.000000		1.000	1.000	83/10/06	83/10/06	
78109	ALLYLCLR TOT WH W	UG/L WATER	K	1	.5000000		.50	.50	83/10/06	83/10/06	
78110	DICLACNI TOT WH W	UG/L WATER	K	1	.5000000		.50	.50	83/10/06	83/10/06	
78121	P-XYLEN+ O-XYLEN	TOT UG/L WATER	K	1	.5000000		.50	.50	83/10/06	83/10/06	
81501	PENTACL ETHANE	TOT UG/L WATER	K	1	2.000000		2.000	2.000	83/10/06	83/10/06	
81552	ACETONE	TOT UG/L WATER	K	1	10.000000		10.000	10.000	83/10/06	83/10/06	
81576	DIETHYL ETHER	TOT UG/L WATER	K	1	1.000000		1.000	1.000	83/10/06	83/10/06	
81595	MTH ETH KETONE	TOT UG/L WATER	K	1	5.000000		5.000	5.000	83/10/06	83/10/06	
81596	MTHISOBU KETONE	TOT UG/L WATER	K	1	1.000000		1.000	1.000	83/10/06	83/10/06	
81607	TETRAHYD FURAN	TOT UG/L WATER	K	1	5.000000		5.000	5.000	83/10/06	83/10/06	
81611	TRICL TRIFLETH	TOT UG/L WATER	K	1	.5000000		.500	.500	83/10/06	83/10/06	
82368	CALCIUM AS CACO3	DIS MG/L WATER		1	180.0000		180.0000	180.0000	78/10/25	78/10/25	
82369	MGSNMIUM AS CACO3	DIS MG/L WATER		1	97.00000		96.9995	96.9995	78/10/25	78/10/25	
84000	GEOLOGIC AGE	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01	
84001	AQUIFER NAME	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01	
84002	CODE GENERAL	REMARKS WATER	TXT	3	TEXT	TEXT	TEXT	TEXT	78/10/25	83/10/06	

STORET RETRIEVAL DATE 87/07/29

PGM=INVENT

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(GWQ0138)

/TYP/A/AMBN/T/WELL

441858093001801 1101820DCAA1 144664  
 44-18 57.9 093 00 18.7 2  
 AQF: PRAIRIE DU CHIEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070639  
 MINOR BASIN: LOWER UPPER MISS RIVER  
 21MINNG 790122 07040002  
 0000 FEET DEPTH

367PRDC

	PARAMETER		MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
00008	LAB IDENT.		WATER		2	125750.0	.0000000	.0000000	125750	125748	83/10/26	83/10/26
00010	WATER TEMP	CENT	WATER		2	9.800000	.0801700	.2831400	10.0	9.6	78/10/26	83/10/26
00089	COLOR PT-CO	UNITS	WATER	K	1	5.000000			5	5	78/10/26	78/10/26
00094	CNDUCTVY FIELD	MICROMHO	WATER		2	474.5000	544.5000	23.33500	491	458	78/10/26	83/10/26
00095	CNDUCTVY AT 25C	MICROMHO	WATER		1	540.0000			540	540	78/10/26	78/10/26
00136	SAMPLE TEMP AT	LAB DEGC	WATER		1	2.500000			2.50000	2.50000	83/10/26	83/10/26
00400	PH	SU	WATER		2	7.200000	.0200040	.1414400	7.30	7.10	78/10/26	83/10/26
00403	LAB PH	SU	WATER		2	7.500000	.0199890	.1413800	7.6	7.4	78/10/26	83/10/26
00410	T ALK CACO3	MG/L	WATER		2	225.0000	50.00000	7.071100	230	220	78/10/26	83/10/26
00425	HCO3 ALK	CACO3	MG/L		2	225.0000	50.00000	7.071100	230	220	78/10/26	83/10/26
00431	T ALK FIELD	MG/L	WATER		2	242.0000	648.0000	25.45600	260	224	78/10/26	83/10/26
00505	RESIDUE TOT VOL	MG/L	WATER	K	1	48.00000			48	48	78/10/26	78/10/26
00615	N02-N TOTAL	MG/L	WATER		1	.0100000			.010	.010	78/10/26	78/10/26
00625	TOT KJEL N	MG/L	WATER		1	.8000000			.800	.800	78/10/26	78/10/26
00630	N02&N03 N-TOTAL	MG/L	WATER	K	1	.0600000			.06	.06	78/10/26	78/10/26
				K	1	.0100000			.01	.01	83/10/26	83/10/26
00665	PHOS-TOT	MG/L	P	TOT	2	.0350000	.0012500	.0353550	.06	.01	78/10/26	83/10/26
00680	T ORG C	C	MG/L		1	.0330000			.033	.033	78/10/26	78/10/26
				K	1	2.000000			2.0	2.0	78/10/26	78/10/26
				K	1	1.000000			1.0	1.0	83/10/26	83/10/26
00900	TOT HARD	CACO3	MG/L	TOT	2	1.500000	.5000000	.7071100	2.0	1.0	78/10/26	83/10/26
00910	CALCITUM	CACO3	MG/L		2	254.0000	512.0000	22.62800	270	238	78/10/26	83/10/26
00920	MGNSIUM	CACO3	MG/L		2	165.0000	450.0000	21.21300	180.0	150.0	78/10/26	83/10/26
00929	SODIUM NA.TOT	MG/L	WATER		2	88.00000	.0000000	.0000000	88.0	88.0	78/10/26	83/10/26
00930	SODIUM NA.DISS	MG/L	WATER		1	6.100000			6.10	6.10	83/10/26	83/10/26
00935	PTSSIUM K.DISS	MG/L	WATER		1	6.300000			6.30	6.30	78/10/26	78/10/26
00937	PTSSIUM K.TOT	MG/L	WATER		1	2.400000			2.40	2.40	78/10/26	78/10/26
00940	CHLORIDE TOTAL	MG/L	WATER	K	1	2.200000			2.20	2.20	83/10/26	83/10/26
				K	1	.6800000			.7	.7	83/10/26	83/10/26
				K	1	.5000000			.5	.5	78/10/26	78/10/26
00945	SULFATE SO4-TOT	MG/L	WATER	TOT	2	.5900000	.0162010	.1272800	.7	.5	78/10/26	83/10/26
00946	SULFATE SO4-DISS	MG/L	WATER		1	45.00000			45	45	83/10/26	83/10/26
00950	FLUORIDE F.DISS	MG/L	WATER		1	52.00000			52.0	52.0	78/10/26	78/10/26
00955	SILICA DISOLVED	MG/L	WATER		1	.3400000			.34	.34	78/10/26	78/10/26
01020	BORON B.DISS	UG/L	WATER		1	8.100000			8.1	8.1	78/10/26	78/10/26
01027	CADMİUM CD.TOT	UG/L	WATER		1	220.0000			220	220	78/10/26	78/10/26
01034	CHROMIUM CR.TOT	UG/L	WATER		1	.1300000			.1	.1	78/10/26	78/10/26
01045	IRON FE,TOT	UG/L	WATER		1	1.000000			1	1	78/10/26	78/10/26
					2	785.0000	198450.0	445.4800	1100	470	78/10/26	83/10/26

STORET RETRIEVAL DATE 87/07/29

PGM=INVENT

PAGE: 8

441858093001801 1101820DCAA1 144664  
 44 18 57.9 093 00 18.7 2  
 AQF: PRAIRIE DU CHIEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070639  
 MINOR BASIN: LOWER UPPER MISS RIVER  
 21MINNG 790122 07040002  
 0000 FEET DEPTH

GWQ0138

367PRDC

/TYP/A/MBNT/WELL

	PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
01051	LEAD PB,TOT	UG/L WATER		1	.8800000			.9	.9	78/10/26	78/10/26
01055	MANGNESE MN	UG/L WATER		2	30.00000	200.0000	14.14200	40.0	20.0	78/10/26	83/10/26
01067	NICKEL NI,TOTAL	UG/L WATER	K	1	1.000000			1	1	78/10/26	78/10/26
01092	ZINC ZN,TOT	UG/L WATER		1	140.0000			140	140	78/10/26	78/10/26
31505	TOT COLI MPN CONF	/100ML WATER	K	2	2.100000	.0200030	.1414300	2	2	78/10/26	83/10/26
31615	FEC COLI MPNECMED	/100ML WATER	K	1	2.000000			2	2	78/10/26	78/10/26
32101	DICLBRMT	TOTUG/L WATER	K	1	.5000000			.5	.5	83/10/26	83/10/26
32102	CARBNTET	TOTUG/L WATER	K	1	.2000000			.2	.2	83/10/26	83/10/26
32104	BROMOFRM WHL-WTR	UG/L WATER	K	1	1.000000			1.0	1.0	83/10/26	83/10/26
32106	CHLRFORM	TOTUG/L WATER	M	1	.2000000			.2	.2	83/10/26	83/10/26
34010	TOLUENE	TOT UG/L WATER	K	1	.5000000			.50	.50	83/10/26	83/10/26
34030	BENZENE	TOT UG/L WATER	K	1	.5000000			.50	.50	83/10/26	83/10/26
34301	CHLOROBNE NZENE	TOTWUG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
34306	CHLORODIBROMOMET	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
34371	ETHYLBEN ZENE	TOTWUG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
34423	METHYLN ECHLORID	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
34475	TETRACHL OROETHYL	TOTWUG/L WATER	K	1	2.000000			2.000	2.000	83/10/26	83/10/26
34480	THALLIUM SEDMG/KG	DRY WGT WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34496	11DICHLO ROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34501	11DICHLO ROETHYLE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34506	111TRICH LOROETHA	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34511	112TRICH LOROETHA	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34516	1122TETR ACHLOROE	TOTWUG/L WATER	K	1	2.000000			2.000	2.000	83/10/26	83/10/26
34531	12DICHLO ROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34536	12DICHLO ROBENZEN	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
34541	12DICHLO ROPROPAN	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34546	12DICHLO ROETHENE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34566	13DICHLO ROBENZEN	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
34571	14DICHLO ROBENZEN	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
34576	2CHLOROE THYLVINY	TOTWUG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
34699	T1,3-DCP TOT WAT	UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34704	C1,3-DCP TOT WAT	UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
39180	TRICHLOR ETHYLENE	TOT UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
70300	RESIDUE DISS-180 C	MG/L WATER		1	290.0000			290	290	78/10/26	78/10/26
71900	MERCURY HG,TOTAL	UG/L WATER		1	.3700000			.4	.4	78/10/26	78/10/26
77093	C-1,2DCE TOTAL	UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
77134	M-XYLENE TOTAL	UG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
77135	O-XYLENE TOTAL	UG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
77166	2,3DCLPR TOTAL	UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26

STORET RETRIEVAL DATE 87/07/29

PGM=INVENT

PAGE: 9  
GWQ0138

441858093001801 1101820DCAA1 144664  
 44.18 57.9 093.00 18.7 2  
 AQF: PRAIRIE DU CHIEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070639  
 MINOR BASIN: LOWER UPPER MISS RIVER  
 21MINNG 790122 07040002  
 0000 FEET DEPTH

/TYP/A/MBNT/WELL

	PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
77168	1,1DCLPR TOTAL	UG/L WATER	K	1	.2000000		.200	.200	.200	83/10/26	83/10/26
77173	1,3DCLPR TOTAL	UG/L WATER	K	1	3.000000		3.000	3.000	3.000	83/10/26	83/10/26
77223	IPROPBNZ TOTAL	UG/L WATER	K	1	.5000000		.500	.500	.500	83/10/26	83/10/26
77443	1,2,3TCP TOTAL	UG/L WATER	K	1	2.000000		2.000	2.000	2.000	83/10/26	83/10/26
77562	1112TCLE TOTAL	UG/L WATER	K	1	.2000000		.200	.200	.200	83/10/26	83/10/26
77596	DBRMETHA TOTAL	UG/L WATER	K	1	1.000000		1.000	1.000	1.000	83/10/26	83/10/26
77651	1,2DBRET TOTAL	UG/L WATER	K	1	1.000000		1.000	1.000	1.000	83/10/26	83/10/26
78109	ALLYLCLR TOT WH W	UG/L WATER	K	1	.5000000		.50	.50	.50	83/10/26	83/10/26
78110	DICLACNI TOT WH W	UG/L WATER	K	1	.5000000		.50	.50	.50	83/10/26	83/10/26
78121	P-XYLEN+ O-XYLEN	TOT UG/L WATER	K	1	.5000000		.50	.50	.50	83/10/26	83/10/26
81501	PENTACL ETHANE	TOT UG/L WATER	K	1	2.000000		2.000	2.000	2.000	83/10/26	83/10/26
81552	ACETONE	TOT UG/L WATER	K	1	10.00000		10.000	10.000	10.000	83/10/26	83/10/26
81576	DIETHYL ETHER	TOT UG/L WATER	K	1	1.000000		1.000	1.000	1.000	83/10/26	83/10/26
81595	MTH ETH KETONE	TOT UG/L WATER	K	1	5.000000		5.000	5.000	5.000	83/10/26	83/10/26
81596	MTHISOBU KETONE	TOT UG/L WATER	K	1	1.000000		1.000	1.000	1.000	83/10/26	83/10/26
81607	TETRAHYD FURAN	TOT UG/L WATER	K	1	5.000000		5.000	5.000	5.000	83/10/26	83/10/26
81611	TRICL TRIFLETH	TOT UG/L WATER	K	1	.5000000		.500	.500	.500	83/10/26	83/10/26
82368	CALCIUM AS CACO3	DIS MG/L WATER		1	170.0000		170.0000	170.0000	170.0000	78/10/26	78/10/26
82369	MGNSIUM AS CACO3	DIS MG/L WATER		1	92.00000		91.9995	91.9995	91.9995	78/10/26	78/10/26
84000	GEOLOGIC AGE	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84001	AQUIFER NAME	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84002	CODE GENERAL	REMARKS WATER	TXT	3	TEXT	TEXT	TEXT	TEXT	TEXT	78/10/26	83/10/26

442216092540601 1111731DCD01 144691

44 22 16.0 092 54 06.0 2

AQF: GALENA DOLOMITE 364GLEN

27049 MINNESOTA GOODHUE

MAJOR BASIN: UPPER MISS 070600

MINOR BASIN: LOWER UPPER MISSISSIPPI

21MINNG 790119 07040002

0000 FEET DEPTH

/TYP/A/MBNT/WELL

	PARAMETER	LAB	IDENT.	NUMBER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
00008	WATER	TEMP	CENT	1	WATER		1	125750.0		125751	125751	83/10/26	83/10/26	
00010	COLOR	PT-CO	UNITS	1	WATER	K	1	9.300000		9.3	9.3	78/10/26	78/10/26	
00088	CNDUCTVY	FIELD	MICROMHO	1	WATER		1	5.000000		5	5	78/10/26	78/10/26	
00094	CNDUCTVY	AT 25C	MICROMHO	1	WATER		1	729.0000		729	729	78/10/26	78/10/26	
00095				1			1	450.0000		450	450	78/10/26	78/10/26	
00400	PH		SU	1	WATER		1	7.500000		7.50	7.50	78/10/26	78/10/26	
00403	LAB	PH	SU	1	WATER		1	7.500000		7.5	7.5	78/10/26	78/10/26	
00410	T ALK	CACO3	MG/L	1	WATER		1	290.0000		290	290	78/10/26	78/10/26	
00425	HCO3 ALK	CACO3	MG/L	1	WATER		1	290.0000		290	290	78/10/26	78/10/26	
00431	T ALK	FIELD	MG/L	1	WATER		1	302.0000		302	302	78/10/26	78/10/26	
00505	RESIDUE	TOT VOL	MG/L	1	WATER		1	160.0000		160	160	78/10/26	78/10/26	
00615	NO2-N	TOTAL	MG/L	1	WATER	K	1	.0100000		.010	.010	78/10/26	78/10/26	
00625	TOT KJEL	N	MG/L	1	WATER		1	.1100000		.110	.110	78/10/26	78/10/26	
00630	NO2&N03	N-TOTAL	MG/L	1	WATER		1	8.500000		8.50	8.50	78/10/26	78/10/26	
00665	PHOS-TOT		MG/L P	1	WATER		1	.0430000		.043	.043	78/10/26	78/10/26	
00680	T ORG C	C	MG/L	1	WATER		1	1.200000		1.2	1.2	78/10/26	78/10/26	
00900	TOT HARD	CACO3	MG/L	1	WATER		1	370.0000		370	370	78/10/26	78/10/26	
00910	CALCIUM	CACO3	MG/L	1	WATER		1	240.0000		240.0	240.0	78/10/26	78/10/26	
00920	MGNSIUM	CACO3	MG/L	1	WATER		1	130.0000		130.0	130.0	78/10/26	78/10/26	
00930	SODIUM	NA,DISS	MG/L	1	WATER		1	4.500000		4.50	4.50	78/10/26	78/10/26	
00935	PTSSUM	K,DISS	MG/L	1	WATER		1	.6600000		.66	.66	78/10/26	78/10/26	
00940	CHLORIDE	TOTAL	MG/L	1	WATER		1	21.00000		21	21	78/10/26	78/10/26	
00946	SULFATE	S04-DISS	MG/L	1	WATER		1	31.00000		31.0	31.0	78/10/26	78/10/26	
00950	FLUORIDE	F,DISS	MG/L	1	WATER		1	.1900000		.19	.19	78/10/26	78/10/26	
00955	SILICA	DISOLVED	MG/L	1	WATER		1	25.00000		25.0	25.0	78/10/26	78/10/26	
01020	BORON	B,DISS	UG/L	1	WATER	K	1	.50.00000		.50	.50	78/10/26	78/10/26	
01027	CADMUM	CD,TOT	UG/L	1	WATER		1	.0300000		.03	.03	78/10/26	78/10/26	
01034	CHROMIUM	CR,TOT	UG/L	1	WATER		1	2.300000		2	2	78/10/26	78/10/26	
01045	IRON	FE,TOT	UG/L	1	WATER	K	1	.50.00000		.50	.50	78/10/26	78/10/26	
01051	LEAD	PB,TOT	UG/L	1	WATER		1	1.000000		1	1	78/10/26	78/10/26	
01055	MANGNESE	MN	UG/L	1	WATER		1	20.00000		20.0	20.0	78/10/26	78/10/26	
01067	NICKEL	NI,TOTAL	UG/L	1	WATER	K	1	1.000000		1	1	78/10/26	78/10/26	
01092	ZINC	ZN,TOT	UG/L	1	WATER		1	24.00000		24	24	78/10/26	78/10/26	
31505	TOT COLI	MPN CONF	/100ML	1	WATER	K	1	2.000000		2	2	78/10/26	78/10/26	
31615	FEC COLI	MPNECMED	/100ML	1	WATER	K	1	2.000000		2	2	78/10/26	78/10/26	
32101	DICLBRMT		TOTUG/L	1	WATER	K	1	.5000000		.5	.5	83/10/26	83/10/26	
32102	CARBNTET		TOTUG/L	1	WATER	K	1	.2000000		.2	.2	83/10/26	83/10/26	
32104	BROMOFORM	WHL-WTR	UG/L	1	WATER	K	1	1.000000		1.0	1.0	83/10/26	83/10/26	
32106	CHLRFORM		TOTUG/L	1	WATER	K	1	.2000000		.2	.2	83/10/26	83/10/26	

STORET RETRIEVAL DATE 87/07/29

PGM=INVENT

PAGE: 11  
GWQ0139

442216092540601 1111731DCD01 144691  
 44 22 16.0 092 54 06.0 2  
 AQF: GALENA DOLomite 364GLEN  
 27049 MINNESOTA GOODHUE  
 MAJOR BASIN: UPPER MISS 070600  
 MINOR BASIN: LOWER UPPER MISSISSIPPI  
 21MINNG 790119 07040002  
 0000 FEET DEPTH

/TYPE/AMNT/WELL

PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
34010 TOLUENE	TOT UG/L WATER	K	1	.5000000			.50	.50	83/10/26	83/10/26
34030 BENZENE	TOT UG/L WATER	K	1	.5000000			.50	.50	83/10/26	83/10/26
34381 CHLOROBENZENE	TOTWUG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
34306 CHLORODIBROMOMET	TOTWUG/L WATER	K	1	1.0000000			1.000	1.000	83/10/26	83/10/26
34371 ETHYLBENZENE	TOTWUG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
34423 METHYLENECHLORIDE	TOTWUG/L WATER	K	1	1.0000000			1.000	1.000	83/10/26	83/10/26
34475 TETRACHLOROETHYL	TOTWUG/L WATER	K	1	2.0000000			2.000	2.000	83/10/26	83/10/26
34480 THALLIUM SEDMG/KG	DRY WGT WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34496 11DICHLOROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34501 11DICHLOROETHYLENE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34506 111TRICHLOROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34511 112TRICHLOROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34516 1122TETRACHLOROETHANE	TOTWUG/L WATER	K	1	2.0000000			2.000	2.000	83/10/26	83/10/26
34531 12DICHLOROETHANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34536 12DICHLOROBENZENE	TOTWUG/L WATER	K	1	1.0000000			1.000	1.000	83/10/26	83/10/26
34541 12DICHLOROPROPANE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34546 12DICHLOROETHENE	TOTWUG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34566 13DICHLOROBENZENE	TOTWUG/L WATER	K	1	1.0000000			1.000	1.000	83/10/26	83/10/26
34571 14DICHLOROBENZENE	TOTWUG/L WATER	K	1	1.0000000			1.000	1.000	83/10/26	83/10/26
34576 2CHLOROETHYL VINYL	TOTWUG/L WATER	K	1	1.0000000			1.000	1.000	83/10/26	83/10/26
34699 T1,3-DCP	TOT WAT UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
34704 C1,3-DCP	TOT WAT UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
39180 TRICHLOROETHYLENE	TOT UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
70300 RESIDUE DISS-180	C MG/L WATER		1	420.0000			420	420	78/10/26	78/10/26
71900 MERCURY HG TOTAL	UG/L WATER		1	.3900000			.4	.4	78/10/26	78/10/26
77093 C-1,2DCE	TOTAL UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
77134 M-XYLENE	TOTAL UG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
77135 O-XYLENE	TOTAL UG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
77166 2,3DCLPR	TOTAL UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
77168 1,1DCLPR	TOTAL UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
77173 1,3DCLPR	TOTAL UG/L WATER	K	1	3.000000			3.000	3.000	83/10/26	83/10/26
77223 IPROPBNZ	TOTAL UG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
77443 1,2,3TCP	TOTAL UG/L WATER	K	1	2.000000			2.000	2.000	83/10/26	83/10/26
77562 1112TCLE	TOTAL UG/L WATER	K	1	.2000000			.200	.200	83/10/26	83/10/26
77596 DBRMETHA	TOTAL UG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
77651 1,2DBRET	TOTAL UG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
78109 ALLYLICLORIDE	TOT WH W UG/L WATER	K	1	.5000000			.50	.50	83/10/26	83/10/26
78110 DICLACNI	TOT WH W UG/L WATER	K	1	.5000000			.50	.50	83/10/26	83/10/26
78121 P-XYLENE	O-XYLEN TOT UG/L WATER	K	1	.5000000			.50	.50	83/10/26	83/10/26

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PGM=INVENT

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GWQ0139

442216092540601 1111731DCD01 144691

44.22 16.0 092 54.06.0 2

AQF: GALENA DOLOMITE 364GLEN

27049 MINNESOTA GOODHUE

MAJOR BASIN: UPPER MISS 070600

MINOR BASIN: LOWER UPPER MISSISSIPPI

21MINNG 790119 07040002

0000 FEET DEPTH

/TYP/A/MBNT/WELL

	PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
81501	PENTACL ETHANE	TOT UG/L WATER	K	1	2.000000			2.000	2.000	83/10/26	83/10/26
81552	ACETONE	TOT UG/L WATER	K	1	10.000000			10.000	10.000	83/10/26	83/10/26
81576	DIETHYL ETHER	TOT UG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
81595	MTH ETH KETONE	TOT UG/L WATER	K	1	5.000000			5.000	5.000	83/10/26	83/10/26
81596	MTHISOBU KETONE	TOT UG/L WATER	K	1	1.000000			1.000	1.000	83/10/26	83/10/26
81607	TETRAHYD FURAN	TOT UG/L WATER	K	1	5.000000			5.000	5.000	83/10/26	83/10/26
81611	TRICL TRIFLETH	TOT UG/L WATER	K	1	.5000000			.500	.500	83/10/26	83/10/26
82368	CALCIUM AS CACO3	DIS MG/L WATER		1	240.0000			240.0000	240.0000	78/10/26	78/10/26
82369	MGNSIUM AS CACO3	DIS MG/L WATER		1	140.0000			140.0000	140.0000	78/10/26	78/10/26
84000	GEOLOGIC AGE	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84001	AQUIFER NAME	CODE WATER	TXT	1	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84002	CODE GENERAL	REMARKS WATER	TXT	2	TEXT	TEXT	TEXT	TEXT	TEXT	78/10/26	83/10/26

4 TOTAL STATIONS PROCESSED

	STA BEG	STA END	# OF OBS	# OF SAMPLE	STA END-PERIOD OF RECD IN YRS
					=0 <.5 <3 >3
<1968	0	0	0	0	0 0 0 0
1968	0	0	0	0	0 0 0 0
1969	0	0	0	0	0 0 0 0
1970	0	0	0	0	0 0 0 0
1971	0	0	0	0	0 0 0 0
1972	0	0	0	0	0 0 0 0
1973	0	0	0	0	0 0 0 0
1974	0	0	0	0	0 0 0 0
1975	0	0	0	0	0 0 0 0
1976	4	0	8	4	0 0 0 0
1977	0	0	0	0	0 0 0 0
1978	0	0	117	3	0 0 0 0
1979	0	0	0	0	0 0 0 0
1980	0	0	0	0	0 0 0 0
1981	0	0	0	0	0 0 0 0
1982	0	0	50	2	0 0 0 0
1983	0	3	194	5	0 0 0 3
1984	0	0	0	0	0 0 0 0
1985	0	0	0	0	0 0 0 0
1986	0	1	66	2	0 0 1 0
1987	0	0	0	0	0 0 0 4
TOTAL	4	4	435	16	0 0 0 4

## 4 TOTAL STATIONS PROCESSED

	PARAMETER	LAB IDENT.	NUMBER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
00005	LAB	IDENT.		WATER		9	125590.0	49152.00	221.7000	125751	125300	82/06/29	86/10/02
00010	WATER	TEMP	CENT	WATER		8	9.550000	.0400390	.2001000	10.0	9.3	78/10/25	86/10/02
00080	COLOR	PT-CO	UNITS	WATER	K	1	5.000000			5	5	78/10/25	78/10/25
					TOT	2	5.000000	.0000000	.0000000	5	5	78/10/28	78/10/26
						3	5.000000	.0000000	.0000000	5	5	78/10/25	78/10/26
00094	CNDUCTVY	FIELD	MICROMHO	WATER		8	524.2500	7480.000	86.48700	729	458	78/10/25	86/10/02
00095	CNDUCTVY	AT 25C	MICROMHO	WATER		4	485.0000	1900.000	43.58900	540	450	78/10/25	82/06/29
00136	SAMPLE TEMP AT	LAB DEGC	WATER			4	3.375000	1.895800	1.376900	5.00000	2.00000	82/06/29	83/10/26
00340	COD	HI LEVEL	MG/L	WATER		1	5.500000			6	6	82/06/29	82/06/29
00400	PH		SU	WATER		7	7.200000	.0633550	.2517000	7.50	6.80	78/10/25	86/10/02
00403	LAB	PH	SU	WATER		7	7.514300	.0148110	.1217000	7.6	7.3	78/10/25	86/10/02
00410	T ALK	CACO3	MG/L	WATER		7	251.4300	514.3000	22.67800	290	220	78/10/25	86/10/02
00425	HCO3 ALK	CACO3	MG/L	WATER		7	251.4300	514.3000	22.67800	290	220	78/10/25	86/10/02
00431	T ALK	FIELD	MG/L	WATER		6	260.0000	657.6000	25.64400	302	224	78/10/25	83/10/26
00505	RESIDUE	TOT VOL	MG/L	WATER	K	4	93.75000	3118.900	55.84700	160	47	78/10/25	82/06/29
00615	NO2-N	TOTAL	MG/L	WATER		3	.0100000	.0000000	.0000000	.010	.010	78/10/25	78/10/26
00625	TOT KJEL	N	MG/L	WATER		4	.4050000	.1055000	.3248100	.800	.110	78/10/25	82/06/29
00630	NO2&NO3	N-TOTAL	MG/L	WATER	K	4	2.152500	17.90800	4.231700	8.50	.01	78/10/25	83/10/06
						3	.0100000	.0000000	.0000000	.01	.01	82/06/29	86/10/02
00665	PHOS-TOT		MG/L P	WATER	TOT	7	1.234300	10.26500	3.203900	8.50	.01	78/10/25	86/10/02
					K	2	.0380000	.0000500	.0070712	.043	.033	78/10/26	78/10/26
					TOT	2	.0030000	.0000080	.0028284	.005	.001	78/10/25	82/06/29
						4	.0205000	.0004276	.0206800	.043	.001	78/10/25	82/06/29
00680	T ORG C	C	MG/L	WATER	K	3	1.466700	.2133400	.4618800	2.0	1.2	78/10/26	86/10/02
					TOT	4	1.000000	.0000000	.0000000	1.0	1.0	78/10/25	83/10/26
						7	1.200000	.1333300	.3651500	2.0	1.0	78/10/25	86/10/02
00900	TOT HARD	CACO3	MG/L	WATER		7	279.7200	1760.700	41.96100	370	238	78/10/25	86/10/02
00910	CALCIUM	CACO3	MG/L	WATER		7	182.8600	790.4900	28.11600	240.0	150.0	78/10/25	86/10/02
00920	MGSNMIUM	CACO3	MG/L	WATER		7	96.42900	226.6300	15.05400	130.0	88.0	78/10/25	86/10/02
00929	SODIUM	NA.TOT	MG/L	WATER		4	4.775000	2.862500	1.691900	6.10	2.30	82/06/29	86/10/02
00930	SODIUM	NA.DISS	MG/L	WATER		3	4.300000	4.440000	2.107100	6.30	2.10	78/10/25	78/10/26
00935	PTSSUM	K.DISS	MG/L	WATER		3	1.353300	.8505400	.9222500	2.40	.66	78/10/25	78/10/26
00937	PTSSUM	K.TOT	MG/L	WATER		4	2.075000	.2758400	.5252000	2.40	1.30	82/06/29	86/10/02
00940	CHLORIDE	TOTAL	MG/L	WATER	K	5	4.716000	82.86700	9.103100	21	.6	78/10/26	86/10/02
						2	.5000000	.0000000	.0000000	.5	.5	78/10/25	78/10/26
						7	3.511400	59.47600	7.712100	21	.5	78/10/25	86/10/02
00945	SULFATE	SO4-TOT	MG/L	WATER		4	36.00000	180.6700	13.44100	45	16	82/06/29	86/10/02
00946	SULFATE	SO4-DISS	MG/L	WATER		3	33.33300	310.3300	17.61600	52.0	17.0	78/10/25	78/10/26
00950	FLUORIDE	F.DISS	MG/L	WATER		3	.2233300	.0108330	.1040800	.34	.14	78/10/25	78/10/26
00951	FLUORIDE	F.TOTAL	MG/L	WATER		1	.3200000			.32	.32	82/06/29	82/06/29
00955	SILICA	DISOLVED	MG/L	WATER		3	15.36700	75.60400	8.695000	25.0	8.1	78/10/25	78/10/26
00956	SILICA	TOTAL	MG/L	WATER		2	9.000000	1.280300	1.131500	9.8	8.2	82/06/29	86/10/02
01002	ARSENIC	AS.TOT	UG/L	WATER	K	1	1.000000			1	1	82/06/29	82/06/29
01007	BARIUM	BA.TOT	UG/L	WATER		1	63.00000			63	63	82/06/29	82/06/29

4 TOTAL STATIONS PROCESSED

01020	BORON	B,DISS	UG/L	WATER	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
							K	.220.0000	.0000000	.0000000	220	220	78/10/26
01022	BORON	B,TOT	UG/L	WATER	TOT	1	120.0000			120	120	82/06/29	82/06/29
						2	.0502500	.0028669	.0535440	.1	.02	78/10/25	82/06/29
						3	106.6700	9633.300	98.15000	220	50	78/10/25	78/10/26
01027	CADMIUM	CD,TOT	UG/L	WATER	TOT	1	1.650000	.8450000	.9192400	.2	.1	78/10/26	78/10/26
						2	.5000000	.0000000	.0000000	.5	.5	78/10/25	82/06/29
						4	.0750000	.7225000	.8500000	.2	.5	78/10/25	82/06/29
01042	COPPER	CU,TOT	UG/L	WATER	TOT	1	1.400000			1	1	82/06/29	82/06/29
						6	452.6800	144840.0	380.5800	1100	6	78/10/25	86/10/02
01045	IRON	FE,TOT	UG/L	WATER	TOT	1	50.00000			50	50	78/10/26	78/10/26
						7	395.1600	143870.0	379.3000	1100	6	78/10/25	86/10/02
						3	1.426700	.7141300	.8450700	2	.9	78/10/25	78/10/26
01051	LEAD	PB,TOT	UG/L	WATER	TOT	1	.2000000			.2	.2	82/06/29	82/06/29
						4	1.120000	.8522700	.9231900	2	.2	78/10/25	82/06/29
						6	26.67200	306.3500	17.50300	50.0	.03	78/10/25	86/10/02
01055	MANGANESE	MN	UG/L	WATER	TOT	1	20.00000			20.0	20.0	82/06/29	82/06/29
						7	25.71900	261.6500	16.17600	50.0	.03	78/10/25	86/10/02
						1	4.600000			5	5	78/10/25	78/10/25
01067	NICKEL	NI,TOTAL	UG/L	WATER	TOT	3	1.000000	.0000000	.0000000	1	1	78/10/25	82/06/29
						4	1.900000	3.240000	1.800000	5	1	78/10/25	82/06/29
						4	118.5000	8835.700	93.99800	240	24	78/10/25	82/06/29
01092	ZINC	ZN,TOT	UG/L	WATER	TOT	1	1.000000			1	1	82/06/29	82/06/29
						1	2.000000			2	2	78/10/25	78/10/25
						5	2.080000	.0120010	.1095500	2	2	78/10/26	83/10/26
31505	TOT COLI	MPN CONF	/100ML	WATER	TOT	1	24000.00			24000	24000	82/06/29	82/06/29
						7	3430.400	82272000	9070.400	24000	2	78/10/25	83/10/26
						4	2.000000	.0000000	.0000000	2	2	78/10/25	82/10/14
31615	FEC COLI	MPNECMED	/100ML	WATER	TOT	1	2.000000			2	2	82/10/14	82/10/14
						4	.5000000	.0000000	.0000000	.5	.5	83/10/06	86/10/02
						4	.2000000	.0000000	.0000000	.2	.2	83/10/06	86/10/02
31679	FECSTREP	MF M-ENT	/100ML	WATER	TOT	4	1.000000	.0000000	.0000000	1.0	1.0	83/10/06	86/10/02
						3	.2000000	.0000000	.0000000	.2	.2	83/10/06	86/10/02
						1	.2000000			.2	.2	83/10/26	83/10/26
32101	DICLBRMT		TOTUG/L	WATER	TOT	4	.5000000			.5	.5	83/10/06	86/10/02
						4	.2000000	.0000000	.0000000	.2	.2	83/10/06	86/10/02
						4	1.000000	.0000000	.0000000	1.0	1.0	83/10/06	86/10/02
32102	CARBTET		TOTUG/L	WATER	TOT	3	.2000000	.0000000	.0000000	.2	.2	83/10/06	86/10/02
						1	.2000000			.2	.2	83/10/26	83/10/26
						4	.2000000	.0000000	.0000000	.2	.2	83/10/06	86/10/02
32104	BROMOFRM	WHL-WTR	UG/L	WATER	TOT	4	.8750000	.0625000	.2500000	1.000	.500	83/10/06	86/10/02
						4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
						4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
32106	CHLRFORM		TOTUG/L	WATER	TOT	4	2.000000	.0000000	.0000000	2.000	2.000	83/10/06	86/10/02
						3	.2000000	.0000000	.0000000	.200	.200	83/10/06	83/10/26
						1	2.000000			.2	.2	83/10/06	86/10/02
32730	PHENOLS	TOTAL	UG/L	WATER	TOT	4	.2000000	.0000000	.0000000	.2	.2	83/10/06	86/10/02
						1	2.000000			.2	.2	82/06/29	82/06/29
						4	.5000000	.0000000	.0000000	.50	.50	83/10/06	86/10/02
34010	TOLUENE		TOTUG/L	WATER	TOT	4	.5000000	.0000000	.0000000	.50	.50	83/10/06	86/10/02
						4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
						4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
34030	BENZENE		TOTUG/L	WATER	TOT	4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
						4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
						4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
34301	CHLOROBENZENE		TOTWUG/L	WATER	TOT	4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
						4	.8750000	.0625000	.2500000	1.000	.500	83/10/06	86/10/02
						4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
34306	CHLORODIBROMOMETHANE		TOTWUG/L	WATER	TOT	4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
						4	2.000000	.0000000	.0000000	2.000	2.000	83/10/06	86/10/02
						3	.2000000	.0000000	.0000000	.200	.200	83/10/06	83/10/26
34371	ETHYLBENZENE		TOTWUG/L	WATER	TOT	4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
						4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
						4	2.000000	.0000000	.0000000	2.000	2.000	83/10/06	86/10/02
34423	METHYLENECHLORIDE		TOTWUG/L	WATER	TOT	4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
						4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
						4	2.000000	.0000000	.0000000	2.000	2.000	83/10/06	86/10/02
34475	TETRACHLOROETHYL		TOTWUG/L	WATER	TOT	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
						3	.2000000	.0000000	.0000000	.200	.200	83/10/06	83/10/26
						1	.2000000			.2	.2	83/10/06	86/10/02
34480	THALLIUM	SEDMG/KG	DRY WGT	WATER	TOT								

4 TOTAL STATIONS PROCESSED

	PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
34488	TRICHLOR OFLUOROM	TOTWUG/L WATER	K	1	.5000000			.500	.500	86/10/02	86/10/02
34496	11DICHLO ROETHANE	TOTWUG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
34501	11DICHLO ROETHYLE	TOTWUG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
34506	111TRICH LOROETHA	TOTWUG/L WATER	K	4	.2750000	.0225000	.1500000	.500	.200	83/10/06	86/10/02
34511	112TRICH LOROETHA	TOTWUG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
34516	1122TETR ACHLOROE	TOTWUG/L WATER	K	4	2.000000	.0000000	.0000000	2.000	2.000	83/10/06	86/10/02
34531	12DICHLO ROETHANE	TOTWUG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
34536	12DICHLO ROBENZEN	TOTWUG/L WATER	K	4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
34541	12DICHLO ROPROPAN	TOTWUG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
34546	12DICHLO ROETHENE	TOTWUG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
34566	13DICHLO ROBENZEN	TOTWUG/L WATER	K	4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
34571	14DICHLO ROBENZEN	TOTWUG/L WATER	K	4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
34576	2CHLOROE THYLVINY	TOTWUG/L WATER	K	3	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	83/10/26
34699	T1,3-DCP TOT WAT	UG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
34704	C1,3-DCP TOT WAT	UG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
39180	TRICHLOR ETHYLENE	TOT UG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
70300	RESIDUE DISS-180 C	MG/L WATER		4	322.5000	4491.700	67.02000	420	270	78/10/25	82/06/29
71900	MERCURY HG,TOTAL	UG/L WATER		3	3933300	.0006337	.0251740	.4	.4	78/10/25	78/10/26
			K	1	1.000000			.1	.1	82/06/29	82/06/29
			TOT	4	.3200000	.0219340	.1481000	.4	.1	78/10/25	82/06/29
74041	WQF SAMPLE	UPDATED WATER		2	865670.0	41943000	6476.400	870205	861125	86/10/02	86/10/02
77093	C-1,2DCE TOTAL	UG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
77134	M-XYLENE TOTAL	UG/L WATER	K	4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
77135	O-XYLENE TOTAL	UG/L WATER	K	3	.5000000	.0000000	.0000000	.500	.500	83/10/06	83/10/26
77166	2,3DCLPR TOTAL	UG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
77168	1,1DCLPR TOTAL	UG/L WATER	K	4	.2000000	.0000000	.0000000	.200	.200	83/10/06	86/10/02
77173	1,3DCLPR TOTAL	UG/L WATER	K	3	3.000000	.0000000	.0000000	3.000	3.000	83/10/06	83/10/26
77223	IPROPBNZ TOTAL	UG/L WATER	K	4	.6250000	.0625000	.2500000	1.000	.500	83/10/06	86/10/02
77443	1,2,3TCP TOTAL	UG/L WATER	K	3	2.000000	.0000000	.0000000	2.000	2.000	83/10/06	83/10/26
77562	1112TCLE TOTAL	UG/L WATER	K	3	.2000000	.0000000	.0000000	.200	.200	83/10/06	83/10/26
77596	DBRMETHA TOTAL	UG/L WATER	K	4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
77651	1,2DBRET TOTAL	UG/L WATER	K	4	.8750000	.0625000	.2500000	1.000	.500	83/10/06	86/10/02
77652	112C122F TOTAL	UG/L WATER	K	1	.2000000			.200	.200	86/10/02	86/10/02
78109	ALLYLCLR TOT WH W	UG/L WATER	K	4	.5000000	.0000000	.0000000	.50	.50	83/10/06	86/10/02
78110	DICLACNI TOT WH W	UG/L WATER	K	3	.5000000	.0000000	.0000000	.50	.50	83/10/06	83/10/26
78121	P-XYLEN+ O-XYLEN	TOT UG/L WATER	K	4	.5000000	.0000000	.0000000	.50	.50	83/10/06	86/10/02
81501	PENTACL ETHANE	TOT UG/L WATER	K	3	2.000000	.0000000	.0000000	2.000	2.000	83/10/06	83/10/26
81552	ACETONE TOT	UG/L WATER	K	4	12.50000	.25.00000	5.00000	20.000	10.000	83/10/06	86/10/02
81576	DIETHYL ETHER	TOT UG/L WATER	K	4	1.000000	.0000000	.0000000	1.000	1.000	83/10/06	86/10/02
81595	MTH ETH KETONE	TOT UG/L WATER	K	4	5.000000	.0000000	.0000000	5.000	5.000	83/10/06	86/10/02
81596	MTHISOBU KETONE	TOT UG/L WATER	K	4	1.250000	.2500000	.5000000	2.000	1.000	83/10/06	86/10/02
81607	TETRAHYD FURAN	TOT UG/L WATER	K	4	6.250000	.6.250000	.2.50000	10.000	5.000	83/10/06	86/10/02
81611	TRICL TRIFLETH	TOT UG/L WATER	K	4	.5000000	.0000000	.0000000	.500	.500	83/10/06	86/10/02
82368	CALCIUM AS CACO3 DIS	MG/L WATER		3	196.6700	1433.400	37.86000	240.0000	170.0000	78/10/25	78/10/26

STORET RETRIEVAL DATE 87/07/29

PGM=INVENT  
GROSS

PAGE: 17

4 TOTAL STATIONS PROCESSED

	PARAMETER	MEDIUM	RMK	NUMBER	MEAN	VARIANCE	STAN DEV	MAXIMUM	MINIMUM	BEG DATE	END DATE
82369	MGNSIUM AS CACO3 DIS	MG/L WATER		3	109.6700	696.3500	26.38800	140.0000	91.9995	78/10/25	78/10/26
84000	GEOLOGIC AGE	CODE WATER	TXT	4	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84001	AQUIFER NAME	CODE WATER	TXT	4	TEXT	TEXT	TEXT	TEXT	TEXT	76/01/01	76/01/01
84002	CODE GENERAL	REMARKS WATER	TXT	12	TEXT	TEXT	TEXT	TEXT	TEXT	78/10/25	86/10/02

STORED RETRIEVAL DATE 85/05/07

440400092451601 1071616CCDC1 215847 GWO0026  
44 03 59.0 092 45 15 8 2  
AQF: ST. PETER SANDSTONE 364STPR  
27039 MINNESOTA DODGE  
MAJOR BASIN: UPPER MISS 070341  
MINOR BASIN: LOWER UPPER MISS RIVER  
21MINNG 780816 07040004  
0000 FEET DEPTH CSN-RSP 0475368-0870954

/TYP&/MUN/INTAKE/AMBNT/WELL

DATE FROM TO	TIME OF DAY	DEPTH FEET	HCO3 CACO3 MG/L	00425 TOT HARD CACO3 MG/L	00900 CHLORIDE TOTAL MG/L	00940 SULFATE SO4-TOT MG/L	00945 SODIUM NA.TOT MG/L	00929 PTSSIU M K.TOT MG/L	00937 FLUORIDE F.TOTAL MG/L	00951 N	00625 TOT KJEL N MG/L	00630 NO2&NO3 N-TOTAL MG/L	00665 PHOS-TOT MG/L P
78/03/23 83/10/96	14 00 13 50		298 280	310	1 0.5K	47	4.00	2.38		0.300K 0.01K	0.02 0.004		

DATE FROM TO	TIME OF DAY	DEPTH FEET	T ORG C C	00680 COD HI LEVEL	00340 PHENOLS MG/L	32730 TOTAL UG/L	31505 TOT COLI MPN CONF /100ML	31615 FEC COLI MPNECMED /100ML	31679 FECSTREP MF M-ENT /100ML	00500 RESIDUE TOTAL MG/L	00505 RESIDUE TOT VOL MG/L	70300 RESIDUE DISS-180 C MG/L	01002 ARSENIC AS, TOT UG/L
78/03/23	14 00			1.0K			2K	2K	9K		120	350	
83/10/06	13 50				1.3			2					
83/10/27	16 20						2K						

DATE FROM TO	TIME OF DAY	DEPTH FEET	01087 BARIUM BA.TOT UG/L	01022 BORON B.TOT UG/L	01027 CADMIUM CD.TOT UG/L	00910 CALCIUM CACO3 MG/L	01034 CHROMIUM CR.TOT UG/L	01042 COPPER CU.TOT UG/L	01045 IRON FE.TOT UG/L	01051 LEAD PB.TOT UG/L	00920 MGSNIUM CACO3 MG/L	01055 MANGNESE MN UG/L
78/03/23 83/10/06	14 00 13 50			0.02		200.0 210.0	1		450 1300	1	110.0 110.0	54.0 50.0

DATE FROM TO	TIME OF DAY	DEPTH FEET	MERCURY HG.TOTAL UG/L	71900	01067	NICKEL NI.TOTAL UG/L	01147	SELENIUM SE.TOT UG/L	00956	SILICA TOTAL MG/L	ZINC ZN.TOT UG/L
78/03/23	14:00			0.1K		2K				45	

STORED RETRIEVAL DATE 85/05/07

/TYP/A/MUN/INTAKE/AMBNT/WELL

440400092451601 1071616CCDC1 215847 GWQ0026  
 44 03 59.0 092 45 15.8 2  
 AQF: ST. PETER SANDSTONE 364STPR  
 27039 MINNESOTA DODGE  
 MAJOR BASIN: UPPER MISS 070341  
 MINOR BASIN: LOWER UPPER MISS RIVER  
 21MINNG 780816 07840004  
 0000 FEET DEPTH CSN-RSP 0475368-0870954

## **DESCRIPTION**

STATION ESTABLISHED BY THE MINNESOTA POLLUTION CONTROL AGENCY IN 1978  
AS PART OF A GROUND WATER QUALITY MONITORING PROGRAM.

NAME: CITY OF MANTORVILLE WELL #1; FIRE STATION

M.ADDRESS: CITY CLERK'S OFFICE, MANTORVILLE, MN 55955

DIRECTIONS: 5TH STREET IN MANTORVILLE. WELL LOCATED JUST SOUTH OF FIRE STATION: SAMPLED FROM INSIDE FAUCET ON E END OF BUILDING.

WELL DATA(OWNERSHIP): CITY      WELL USE: WITHDRAW      YEAR CONSTR: 1959

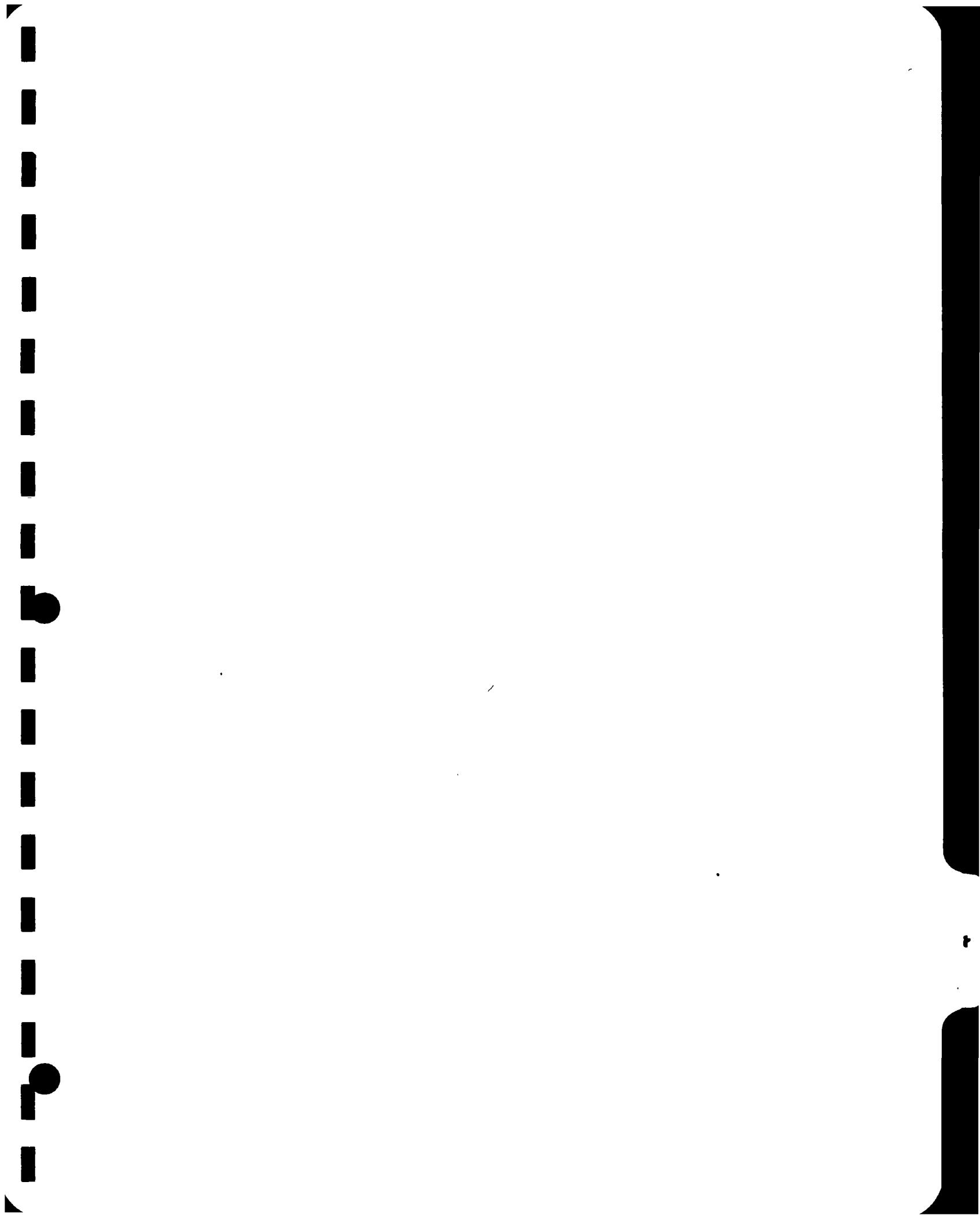
DEPTH FT: 181 LAND SURF DATUM FT: 1125 WATER USE: PS STANDBY, FIRE

CASING(FT CASED): 155 DIAM INCHES: 8 MATERIAL: STEEL

OBSERVATIONS: SEE GROUND WATER MONITORING PHOTO LOG FOR THE EXACT SITE

LOCATION. PARTIAL WELL LOG AVAILABLE. SLIGHT IRON ODOR B31006.

QUAD: DODGE CENTER 7.5'





POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 1 - SITE LOCATION AND INSPECTION INFORMATION

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER
MND	981961873

II. SITE NAME AND LOCATION

01 SITE NAME (Legal, common, or descriptive name of site)	02 STREET, ROUTE NO., OR SPECIFIC LOCATION IDENTIFIER				
HADER GROUND WATER CONTAMINATION	NE 1/4 SEC. 5, T110N, R16W				
03 CITY	04 STATE	05 ZIP CODE	06 COUNTY	07 COUNTY CODE	08 CONG DIST
MINNEOLA TOWNSHIP	MN	55027	GOODHUE	049	01
09 COORDINATES LATITUDE <b>44° 22' 30.0"</b>	LONGITUDE <b>092° 45' 00.0"</b>	10 TYPE OF OWNERSHIP (Check one) <input type="checkbox"/> A PRIVATE <input type="checkbox"/> B FEDERAL <input type="checkbox"/> C STATE <input type="checkbox"/> D COUNTY <input type="checkbox"/> E MUNICIPAL <input checked="" type="checkbox"/> F OTHER <i>private property in rural/agricultural area</i> <input type="checkbox"/> G UNKNOWN			

III. INSPECTION INFORMATION

01 DATE OF INSPECTION <b>6/1/87</b> MONTH DAY YEAR	02 SITE STATUS <b>N/A</b>	03 YEARS OF OPERATION BEGNING YEAR _____ ENDING YEAR _____	<input checked="" type="checkbox"/> UNKNOWN
04 AGENCY PERFORMING INSPECTION (Check all that apply) <input type="checkbox"/> A EPA <input type="checkbox"/> B EPA CONTRACTOR _____ <input checked="" type="checkbox"/> E. STATE <input type="checkbox"/> F. STATE CONTRACTOR _____			
		(Name of firm) _____	
		(Name of firm) _____	
05 CHIEF INSPECTOR <b>BECKY LOFGREN</b>	06 TITLE <b>PROJECT MANAGER</b>	07 ORGANIZATION <b>MPCA</b>	08 TELEPHONE NO <b>(612) 296-7391</b>
09 OTHER INSPECTORS <b>SHAWN LUOTSINOJA</b>	10 TITLE <b>ASSIS. PROJECT MANAGER</b>	11 ORGANIZATION <b>MPCA</b>	12 TELEPHONE NO <b>(612) 296-7783</b>
<b>SUSAN PRICE</b>	<b>HYDROLOGIST</b>	<b>MPCA</b>	<b>(612) 297-1796</b>
			( )
			( )
			( )
13 SITE REPRESENTATIVES INTERVIEWED <b>non responsive</b>	14 TITLE	15 ADDRESS	16 TELEPHONE NO
			( )
			( )
			( )
			( )
			( )

17 ACCESS GAINED BY (Check one) <input checked="" type="checkbox"/> PERMISSION <input type="checkbox"/> WARRANT	18 TIME OF INSPECTION <b>3-29-88 to 3-31-88 8:00 AM - 5:00 PM</b>	19 WEATHER CONDITIONS <b>SUNNY AND COOL</b>
--	--	--

IV. INFORMATION AVAILABLE FROM

01 CONTACT <b>SUSAN PRICE</b>	02 OF (Agency/Organization) <b>MN POLLUTION CONTROL AGENCY (MPCA)</b>	03 TELEPHONE NO <b>(612) 297-1796</b>
04 PERSON RESPONSIBLE FOR SITE INSPECTION FORM <b>BECKY LOFGREN</b>	05 AGENCY <b>MPCA</b>	06 ORGANIZATION 07 TELEPHONE NO <b>(612) 296-7391</b>
		08 DATE _____ MONTH DAY YEAR



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 2 - WASTE INFORMATION

01 STATE	02 SITE NUMBER
MND	981961873

II. WASTE STATES, QUANTITIES, AND CHARACTERISTICS

01 PHYSICAL STATES (Check all that apply)		02 WASTE QUANTITY AT SITE <small>(MEASURES OF WASTE QUANTITY MUST BE INDEPENDENT)</small>		03 WASTE CHARACTERISTICS (Check all that apply)			
<input type="checkbox"/> A SOLID	<input type="checkbox"/> E SLURRY	TONS _____	<input checked="" type="checkbox"/> A TOXIC	<input type="checkbox"/> E SOLUBLE	<input type="checkbox"/> I HIGHLY VOLATILE		
<input type="checkbox"/> B POWDER/FINES	<input checked="" type="checkbox"/> F LIQUID	CUBIC YARDS _____	<input checked="" type="checkbox"/> B CORROSIVE	<input type="checkbox"/> F INFECTIOUS	<input type="checkbox"/> J EXPLOSIVE		
<input type="checkbox"/> C SLUDGE	<input type="checkbox"/> G GAS	NO OF DRUMS _____	<input checked="" type="checkbox"/> C RADIOACTIVE	<input type="checkbox"/> G FLAMMABLE	<input type="checkbox"/> K REACTIVE		
<input type="checkbox"/> D OTHER _____				<input checked="" type="checkbox"/> D PERSISTENT	<input type="checkbox"/> H IGNITABLE	<input type="checkbox"/> L INCOMPATIBLE	
						<input type="checkbox"/> M NOT APPLICABLE	

III. WASTE TYPE

CATEGORY	SUBSTANCE NAME	01 GROSS AMOUNT	02 UNIT OF MEASURE	03 COMMENTS
SLU	SLUDGE			
OLW	OILY WASTE			
SOL	SOLVENTS			
PSD	PESTICIDES			
OCC	OTHER ORGANIC CHEMICALS			
IOC	INORGANIC CHEMICALS			
ACD	ACIDS			
BAS	BASES			
(MES)	HEAVY METALS	UNKNOWN		

IV. HAZARDOUS SUBSTANCES (See ADDENDUM 10: MOST COMMON CAS NUMBERS)

01 CATEGORY	02 SUBSTANCE NAME	03 CAS NUMBER	04 STORAGE/DISPOSAL METHOD	05 CONCENTRATION	06 MEASURE OF CONCENTRATION
MES	Aluminum	7429905		32100	ug/l (ppb)
	Arsenic	7440382		8.4	
	Barium	7440393		91.5	
	Calcium	7440702		167000	
	Chromium	7440473		11.0	
	Cobalt	—		113	
	Copper	7440508		27.4	
	Iron	NA		236000	
	Lead	7439921		19.9	
	Magnesium	7439954		47400	
	Manganese	7439965		1670	
	Nickel	7440020		434	
	Potassium	7440097		56400	
	Sodium	7440235		16680	
	Vanadium	7440622		25.8	
	Zinc	7440600		4120	
	Sulfate	—		890	

V. FEEDSTOCKS (See ADDENDUM 10: CAS NUMBERS)

CATEGORY	D1 FEEDSTOCK NAME	D2 CAS NUMBER	CATEGORY	D1 FEEDSTOCK NAME	D2 CAS NUMBER
FDS	N/A		FDS		
FDS			FDS		
FDS			FDS		
FDS			FDS		

VI. SOURCES OF INFORMATION (See ADDENDUM 10: SITES THAT SUPPORTED ANALYSIS RESULTS)

MPCA Hader GW Contamination Files.

POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT

## PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS

L IDENTIFICATION	
O1 STATE	O2 SITE NUMBER
MND	981961873

## II. HAZARDOUS CONDITIONS AND INCIDENTS

O1  A. GROUNDWATER CONTAMINATION

O3 POPULATION POTENTIALLY AFFECTED

473

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

The elevated concentrations of metals in the residential well water appears to be a naturally occurring problem. The geology of the area and the possibility of iron bacteria in the well could be impacting the quality of the water. No contaminant source has been identified through the site inspection process.

O1  B SURFACE WATER CONTAMINATION

O3 POPULATION POTENTIALLY AFFECTED

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

Numerous small streams and creeks flow through the Site and are used primarily by livestock. However, no source for a release of contaminants has been identified, therefore, surface water contamination is not of concern.

O1  C CONTAMINATION OF AIR

O3 POPULATION POTENTIALLY AFFECTED

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

There is no documentation of air contamination and there is no potential for air contamination at the Site.

O1  D FIRE/EXPLOSIVE CONDITIONS

O3 POPULATION POTENTIALLY AFFECTED

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

Fire and explosive conditions do not exist at this site.

O1  E. DIRECT CONTACT

O3 POPULATION POTENTIALLY AFFECTED

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

No source of the contamination has been identified. The elevated levels of metals appear to be a naturally occurring problem, therefore direct contact is not possible.

O1  F CONTAMINATION OF SOIL

O3 AREA POTENTIALLY AFFECTED

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

No soil samples were taken because of the large area of the Site (approximately 24 square miles) and no source of the contamination could be identified.

O1  G DRINKING WATER CONTAMINATION

O3 POPULATION POTENTIALLY AFFECTED

473

O2  OBSERVED (DATE 3/29-3/1/88)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

Numerous residential wells were sampled during the site inspection. Water from the residential wells show elevated levels of metals - nickel in particular. The ground water in the area is classified as a "calcium-magnesium bicarbonate type". High metal concentrations may be due to the geology of the area and possibly iron bacteria.

O1  H WORKER EXPOSURE/INJURY

O3 WORKERS POTENTIALLY AFFECTED

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

Contamination problem is with residential wells. No workers are "on-site".

O1  I POPULATION EXPOSURE/INJURY

O3 POPULATION POTENTIALLY AFFECTED

O2  OBSERVED (DATE \_\_\_\_\_)

O4 NARRATIVE DESCRIPTION

 POTENTIAL ALLEGED

The only possibility of exposure is through the drinking water.

POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT

## PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER
MND	981961873

## II. HAZARDOUS CONDITIONS AND INCIDENTS (continued)

D1  J DAMAGE TO FLORA  
D4 NARRATIVE DESCRIPTIOND2  OBSERVED (DATE \_\_\_\_\_) D POTENTIAL D ALLEGED

Damage to flora has not been documented and no damage was seen during the site inspection.

D1  K DAMAGE TO FAUNA  
D4 NARRATIVE DESCRIPTION (include names of species)D2  OBSERVED (DATE \_\_\_\_\_) D POTENTIAL D ALLEGED

Local farmers allege that livestock have died from drinking "bad" water. There is no documented proof of this claim.

D1  L CONTAMINATION OF FOOD CHAIN  
D4 NARRATIVE DESCRIPTIOND2  OBSERVED (DATE \_\_\_\_\_) D POTENTIAL D ALLEGED

SEE K. ABOVE

D1  M UNSTABLE CONTAINMENT OF WASTES  
(Soils Runoff Standing Pools Leaking Drums)D2  OBSERVED (DATE \_\_\_\_\_) D POTENTIAL D ALLEGED

D3 POPULATION POTENTIALLY AFFECTED \_\_\_\_\_

D4 NARRATIVE DESCRIPTION

There is no evidence of hazardous waste on-site.

D1  N DAMAGE TO OFFSITE PROPERTY  
D4 NARRATIVE DESCRIPTIOND2  OBSERVED (DATE \_\_\_\_\_) D POTENTIAL D ALLEGED

Other offsite residential wells may have the same problem with elevated metal levels.

D1  O CONTAMINATION OF SEWERS STORM DRAINS WWTPs  
D4 NARRATIVE DESCRIPTIOND2  OBSERVED (DATE \_\_\_\_\_) D POTENTIAL D ALLEGED

Sewer and storm drains are not in the area of concern. This is a rural area with septic systems.

D1  P ILLEGAL/UNAUTHORIZED DUMPING  
D4 NARRATIVE DESCRIPTIOND2  OBSERVED (DATE \_\_\_\_\_) D POTENTIAL D ALLEGED

An unpermitted dump had been operating from the early 1970s through 1980 in an abandoned quarry approximately 1½ miles northwest of the Site. This dump is being investigated through a site inspection on the Hader Dump.

D5 DESCRIPTION OF ANY OTHER KNOWN POTENTIAL OR ALLEGED HAZARDS

III. TOTAL POPULATION POTENTIALLY AFFECTED \_\_\_\_\_

IV. COMMENTS  
\_\_\_\_\_  
\_\_\_\_\_

V. SOURCES OF INFORMATION (Cite specific references e.g. State and County Health Departments)

MPCA Hader GW Contamination File



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION  
PART 4 - PERMIT AND DESCRIPTIVE INFORMATION

I. IDENTIFICATION

01 STATE	02 SITE NUMBER
MN	D981961873

II. PERMIT INFORMATION

01 TYPE OF PERMIT ISSUED  
(Check all that apply)

- A NPDES
- B UIC
- C AIR
- D RCRA
- E RCRA INTERIM STATUS
- F SPCC PLAN
- G STATE (Specify)
- H. LOCAL (Specify)
- I OTHER (Specify)
- J NONE

02 PERMIT NUMBER

03 DATE ISSUED

04 EXPIRATION DATE

05 COMMENTS

III. SITE DESCRIPTION

01 STORAGE/DISPOSAL (Check all that apply)

02 AMOUNT

N/A

03 UNIT OF MEASURE

- A. SURFACE IMPOUNDMENT
- B. PILES
- C. DRUMS, ABOVE GROUND
- D. TANK, ABOVE GROUND
- E. TANK, BELOW GROUND
- F. LANDFILL
- G. LANDFARM
- H. OPEN DUMP
- I. OTHER \_\_\_\_\_  
(Specify)

04 TREATMENT (Check all that apply)

N/A

- A. INCINERATION
- B. UNDERGROUND INJECTION
- C. CHEMICAL/PHYSICAL
- D. BIOLOGICAL
- E. WASTE OIL PROCESSING
- F. SOLVENT RECOVERY
- G. OTHER RECYCLING/RECOVERY
- H. OTHER \_\_\_\_\_  
(Specify)

05 OTHER

- A. BUILDINGS ON SITE

Several Farmsteads

06 AREA OF SITE

15.360 (Acres)

07 COMMENTS

IV. CONTAINMENT N/A

01 CONTAINMENT OF WASTES (Check one)

- A. ADEQUATE, SECURE
- B. MODERATE
- C. INADEQUATE, POOR
- D. INSECURE, UNSOUND, DANGEROUS

02 DESCRIPTION OF DRUMS, DIKING, LINERS, BARRIERS, ETC

N/A

V. ACCESSIBILITY

01 WASTE EASILY ACCESSIBLE  YES  NO

02 COMMENTS

VI. SOURCES OF INFORMATION (Give specific references e.g. State files, sample analysis, reports)

MPCA Hader G W Contamination File



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 5 - WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER

MN D981961873

II. DRINKING WATER SUPPLY

01 TYPE OF DRINKING SUPPLY  
(Check as applicable)

SURFACE	WELL
COMMUNITY	A <input type="checkbox"/> B <input type="checkbox"/>
NON-COMMUNITY	C <input type="checkbox"/> D <input checked="" type="checkbox"/>

02 STATUS

ENDANGERED	AFFECTED	MONITORED
A <input type="checkbox"/>	B. <input type="checkbox"/>	C. <input type="checkbox"/>
D <input type="checkbox"/>	E <input checked="" type="checkbox"/>	F. <input type="checkbox"/>

03 DISTANCE TO SITE

A \_\_\_\_\_ (mi)  
B \_\_\_\_\_ (mi)

III. GROUNDWATER

01 GROUNDWATER USE IN VICINITY (Check one)

A ONLY SOURCE FOR DRINKING     B DRINKING  
(Other sources available)  
COMMERCIAL, INDUSTRIAL, IRRIGATION  
(No other water sources available)

02 POPULATION SERVED BY GROUND WATER 3.8 x 130 = 494

03 DISTANCE TO NEAREST DRINKING WATER WELL 0 (mi)

04 DEPTH TO GROUNDWATER

16 (ft)

05 DIRECTION OF GROUNDWATER FLOW

North

06 DEPTH TO AQUIFER OF CONCERN

16 (ft)

07 POTENTIAL YIELD OF AQUIFER

150 (gpd)

08 SOLE SOURCE AQUIFER

YES  NO

09 DESCRIPTION OF WELLS (Including usage, depth and location relative to population and buildings)

All wells are open-hole to bedrock (Prairie du Chien and Jordan) with the exception of 1 surficial well.

10 RECHARGE AREA

YES COMMENTS Recharge area for Cannon River Watershed  
 NO

11 DISCHARGE AREA

YES COMMENTS  
 NO

IV. SURFACE WATER

01 SURFACE WATER USE (Check one)

A. RESERVOIR, RECREATION DRINKING WATER SOURCE     B. IRRIGATION, ECONOMICALLY IMPORTANT RESOURCES     C. COMMERCIAL, INDUSTRIAL     D. NOT CURRENTLY USED  
Occasionally used by livestock for drinking

02 AFFECTED/POTENTIALLY AFFECTED BODIES OF WATER

NAME:

Belle Creek  
Zumbro River

AFFECTED

DISTANCE TO SITE

1.5 (mi)

4.0 (mi)

             (mi)

V. DEMOGRAPHIC AND PROPERTY INFORMATION

01 TOTAL POPULATION WITHIN

ONE (1) MILE OF SITE  
A 141  
NO OF PERSONS

TWO (2) MILES OF SITE  
B 211  
NO OF PERSONS

THREE (3) MILES OF SITE  
C 494  
NO OF PERSONS

02 DISTANCE TO NEAREST POPULATION

0 (mi)

03 NUMBER OF BUILDINGS WITHIN TWO (2) MILES OF SITE

55

04 DISTANCE TO NEAREST OFF-SITE BUILDING

0 (mi)

05 POPULATION WITHIN VICINITY OF SITE (Provide narrative description of nature of population within vicinity of site e.g. rural, village, densely populated urban area)

Minneota and Belle Creek townships are sparsely populated areas in Goodhue county. The land is used primarily for agricultural purposes.



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 5 - WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA

I. IDENTIFICATION

01 STATE MN 02 SITE NUMBER D981961873

VI. ENVIRONMENTAL INFORMATION

01 PERMEABILITY OF UNSATURATED ZONE (Check one)

A  $10^{-6} - 10^{-8}$  cm/sec    B  $10^{-4} - 10^{-6}$  cm/sec    C  $10^{-4} - 10^{-3}$  cm/sec    D GREATER THAN  $10^{-3}$  cm/sec

02 PERMEABILITY OF BEDROCK (Check one)

A IMPERMEABLE (Less than  $10^{-6}$  cm/sec)    B RELATIVELY IMPERMEABLE ( $10^{-4} - 10^{-6}$  cm/sec)    C RELATIVELY PERMEABLE ( $10^{-2} - 10^{-4}$  cm/sec)    D VERY PERMEABLE (Greater than  $10^{-2}$  cm/sec)

03 DEPTH TO BEDROCK

04 DEPTH OF CONTAMINATED SOIL ZONE

05 SOIL pH

35 (ft)

N/A (ft)

N/A

06 NET PRECIPITATION

07 ONE YEAR 24 HOUR RAINFALL

08 SLOPE SITE SLOPE

DIRECTION OF SITE SLOPE

TERRAIN AVERAGE SLOPE

-3 (in)

2.45 (in)

5 %

VARIES DUE TO

5-7 %

LARGE AREA

09 FLOOD POTENTIAL

10

SITE IS ON BARRIER ISLAND, COASTAL HIGH HAZARD AREA, RIVERINE FLOODWAY

11 DISTANCE TO WETLANDS (5 acre minimum)

12 DISTANCE TO CRITICAL HABITAT (of endangered species)

ESTUARINE

OTHER

BELLE CREEK

N/A (mi)

A N/A (mi)

B 0.16 (mi)

ENDANGERED SPECIES \_\_\_\_\_

13 LAND USE IN VICINITY

DISTANCE TO

COMMERCIAL/INDUSTRIAL

RESIDENTIAL AREAS, NATIONAL/STATE PARKS,  
FORESTS, OR WILDLIFE RESERVES

AGRICULTURAL LANDS  
PRIME AG LAND AG LAND

A N/A (mi)

B 0 (mi)

C 0 (mi) D 24 Sq. (mi)

14 DESCRIPTION OF SITE IN RELATION TO SURROUNDING TOPOGRAPHY

Site is hummocky farm land over 24 square miles in area.

VII. SOURCES OF INFORMATION (Check specific references e.g., state files, sample analysis, reports)

MPCA Hader Groundwater Contamination File

Water Resources of the Cannon River Watershed,  
Southeastern Minnesota, Anderson, H.W. et.al. USGS 1971



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 6 - SAMPLE AND FIELD INFORMATION

I. IDENTIFICATION  
01 STATE 02 SITE NUMBER  
MN A981961873

II. SAMPLES TAKEN

SAMPLE TYPE	01 NUMBER OF SAMPLES TAKEN	02 SAMPLES SENT TO	03 ESTIMATED DATE RESULTS AVAILABLE
GROUNDWATER	10	S-Cubed (Organic), ENSECO/Rocky Mtn. Analy. (Inorganic), Post Buckley Schuh, Jernigan (SAS)	
SURFACE WATER	0		
WASTE			
AIR			
RUNOFF			
SPILL			
SOIL			
VEGETATION			
OTHER			

III. FIELD MEASUREMENTS TAKEN

01 TYPE	02 COMMENTS
pH	used to stabilize well prior to sampling
Temperature	
Specific Conductivity	

IV. PHOTOGRAPHS AND MAPS

01 TYPE <input checked="" type="checkbox"/> GROUND <input type="checkbox"/> AERIAL	02 IN CUSTODY OF <u>MN POLLUTION CONTROL AGENCY - SITE ASSESSMENT UNIT</u> <small>(Name of organization or individual)</small>
MAPS <input checked="" type="checkbox"/> YES <input type="checkbox"/> NO	04 LOCATION OF MAPS <u>MN POLLUTION CONTROL AGENCY</u>

OTHER FIELD DATA COLLECTED (Provide narrative description)


V. SOURCES OF INFORMATION (cite specific references e.g. state files, sample analysis reports)

<u>MPCA Hader GW Contamination File</u>



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 7 - OWNER INFORMATION

I. IDENTIFICATION  
01 STATE | 02 SITE NUMBER  
MN | 0981961873

II. CURRENT OWNER(S)			PARENT COMPANY (if applicable)			
01 NAME	02 D+B NUMBER	08 NAME	09 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD #, etc.)		04 SIC CODE	10 STREET ADDRESS (P O Box, RFD #, etc.)			11 SIC CODE
05 CITY	06 STATE	07 ZIP CODE	12 CITY	13 STATE	14 ZIP CODE	
01 NAME	02 D+B NUMBER	08 NAME	09 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD #, etc.)		04 SIC CODE	10 STREET ADDRESS (P O Box, RFD #, etc.)			11 SIC CODE
05 CITY	06 STATE	07 ZIP CODE	12 CITY	13 STATE	14 ZIP CODE	
01 NAME	02 D+B NUMBER	08 NAME	09 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD #, etc.)		04 SIC CODE	10 STREET ADDRESS (P O Box, RFD #, etc.)			11 SIC CODE
05 CITY	06 STATE	07 ZIP CODE	12 CITY	13 STATE	14 ZIP CODE	
01 NAME	02 D+B NUMBER	08 NAME	09 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD #, etc.)		04 SIC CODE	10 STREET ADDRESS (P O Box, RFD #, etc.)			11 SIC CODE
05 CITY	06 STATE	07 ZIP CODE	12 CITY	13 STATE	14 ZIP CODE	
III. PREVIOUS OWNER(S) (List most recent first)			IV. REALTY OWNER(S) (if applicable, list most recent first)			
01 NAME	02 D+B NUMBER	01 NAME	02 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD #, etc.)		04 SIC CODE	03 STREET ADDRESS (P O Box, RFD #, etc.)			04 SIC CODE
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE	
01 NAME	02 D+B NUMBER	01 NAME	02 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD #, etc.)		04 SIC CODE	03 STREET ADDRESS (P O Box, RFD #, etc.)			04 SIC CODE
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE	
01 NAME	02 D+B NUMBER	01 NAME	02 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD #, etc.)		04 SIC CODE	03 STREET ADDRESS (P O Box, RFD #, etc.)			04 SIC CODE
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE	
V. SOURCES OF INFORMATION (Cite specific references, e.g., site files, sample analysis reports)						



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 8 - OPERATOR INFORMATION

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER
MN	D981961873

II. CURRENT OPERATOR (Provide # different from owner)			OPERATOR'S PARENT COMPANY (If applicable)			
01 NAME <i>N/A</i>	02 D+B NUMBER	10 NAME <i>N/A</i>	11 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD#, etc.)		04 SIC CODE	12 STREET ADDRESS (P O Box, RFD#, etc.)		13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER					
III. PREVIOUS OPERATOR(S) (List most recent first; provide only # different from owner)			PREVIOUS OPERATORS' PARENT COMPANIES (If applicable)			
01 NAME <i>N/A</i>	02 D+B NUMBER	10 NAME <i>N/A</i>	11 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD#, etc.)		04 SIC CODE	12 STREET ADDRESS (P O Box, RFD#, etc.)		13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD					
01 NAME	02 D+B NUMBER	10 NAME	11 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD#, etc.)		04 SIC CODE	12 STREET ADDRESS (P O Box, RFD#, etc.)		13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD					
01 NAME	02 D+B NUMBER	10 NAME	11 D+B NUMBER			
03 STREET ADDRESS (P O Box, RFD#, etc.)		04 SIC CODE	12 STREET ADDRESS (P O Box, RFD#, etc.)		13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD					

IV. SOURCES OF INFORMATION (One specific references e.g. state IDES sample analysis reports)

*MPCA Site Inspection*



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 9 - GENERATOR/TRANSPORTER INFORMATION

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER
MN	D981961873

II. ON-SITE GENERATOR

01 NAME <b>N/A</b>	02 D+B NUMBER		
03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE		
05 CITY	06 STATE		

III. OFF-SITE GENERATOR(S)

01 NAME <b>N/A</b>	02 D+B NUMBER	01 NAME	02 D+B NUMBER		
03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE	03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE
01 NAME	02 D+B NUMBER	01 NAME	02 D+B NUMBER		
03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE	03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE

IV. TRANSPORTER(S)

01 NAME <b>N/A</b>	02 D+B NUMBER	01 NAME	02 D+B NUMBER		
03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE	03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE
01 NAME	02 D+B NUMBER	01 NAME	02 D+B NUMBER		
03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE	03 STREET ADDRESS (P O Box, RFD#, etc)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE

V. SOURCES OF INFORMATION (List specific references, e.g., state files, sample analysis, reports)

**MPCA Site Inspection**



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 10 - PAST RESPONSE ACTIVITIES

I. IDENTIFICATION  
01 STATE MN  
02 SITE NUMBER 0981961873

II. PAST RESPONSE ACTIVITIES

01 <input checked="" type="checkbox"/> A. WATER SUPPLY CLOSED 04 DESCRIPTION Discontinued use of their water for drinking and food preparation was advised for 3 residents due to elevated lead and nickel levels. The drinking water advisories still allowed the water to be used for bathing, dishwashing, etc.	02 DATE 9-2-87	03 AGENCY MN DEPT. OF HEALTH
01 <input checked="" type="checkbox"/> B. TEMPORARY WATER SUPPLY PROVIDED 04 DESCRIPTION Bottled water was supplied to 3 residents in the area due to Minnesota Department of Health analytical results showing elevated levels of lead and nickel above recommended drinking water limits.	02 DATE 7-28-87	03 AGENCY MPCA
01 <input type="checkbox"/> C. PERMANENT WATER SUPPLY PROVIDED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> D. SPILLED MATERIAL REMOVED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> E. CONTAMINATED SOIL REMOVED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> F. WASTE REPACKAGED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> G. WASTE DISPOSED ELSEWHERE 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> H. ON SITE BURIAL 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> I. IN SITU CHEMICAL TREATMENT 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> J. IN SITU BIOLOGICAL TREATMENT 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> K. IN SITU PHYSICAL TREATMENT 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> L. ENCAPSULATION 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> M. EMERGENCY WASTE TREATMENT 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> N. CUTOFF WALLS 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> O. EMERGENCY DIKING/SURFACE WATER DIVERSION 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> P. CUTOFF TRENCHES/SUMP 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		
01 <input type="checkbox"/> Q. SUBSURFACE CUTOFF WALL 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
N/A		



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 10 - PAST RESPONSE ACTIVITIES

I. IDENTIFICATION  
01 STATE MN 02 SITE NUMBER D981961873

II PAST RESPONSE ACTIVITIES (Continued)

01 <input type="checkbox"/> R. BARRIER WALLS CONSTRUCTED 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> S. CAPPING/COVERING 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> T. BULK TANKAGE REPAIRED 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> U. GROUT CURTAIN CONSTRUCTED 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> V. BOTTOM SEALED 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> W. GAS CONTROL 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> X. FIRE CONTROL 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> Y. LEACHATE TREATMENT 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> Z. AREA EVACUATED 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> 1. ACCESS TO SITE RESTRICTED 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> 2. POPULATION RELOCATED 04 DESCRIPTION	N/A	02 DATE	03 AGENCY
01 <input type="checkbox"/> 3. OTHER REMEDIAL ACTIVITIES 04 DESCRIPTION	N/A	02 DATE	03 AGENCY

III. SOURCES OF INFORMATION (List specific references, e.g., state files, sample analysis reports)

MPCA Site Inspection



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 11 - ENFORCEMENT INFORMATION

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER
MN	D981961B73

II. ENFORCEMENT INFORMATION

01 PAST REGULATORY/ENFORCEMENT ACTION  YES  NO

02 DESCRIPTION OF FEDERAL, STATE, LOCAL REGULATORY/ENFORCEMENT ACTION

N/A

III. SOURCES OF INFORMATION (List specific references e.g. state files, sample analysis reports)

MPCA GWSW Files

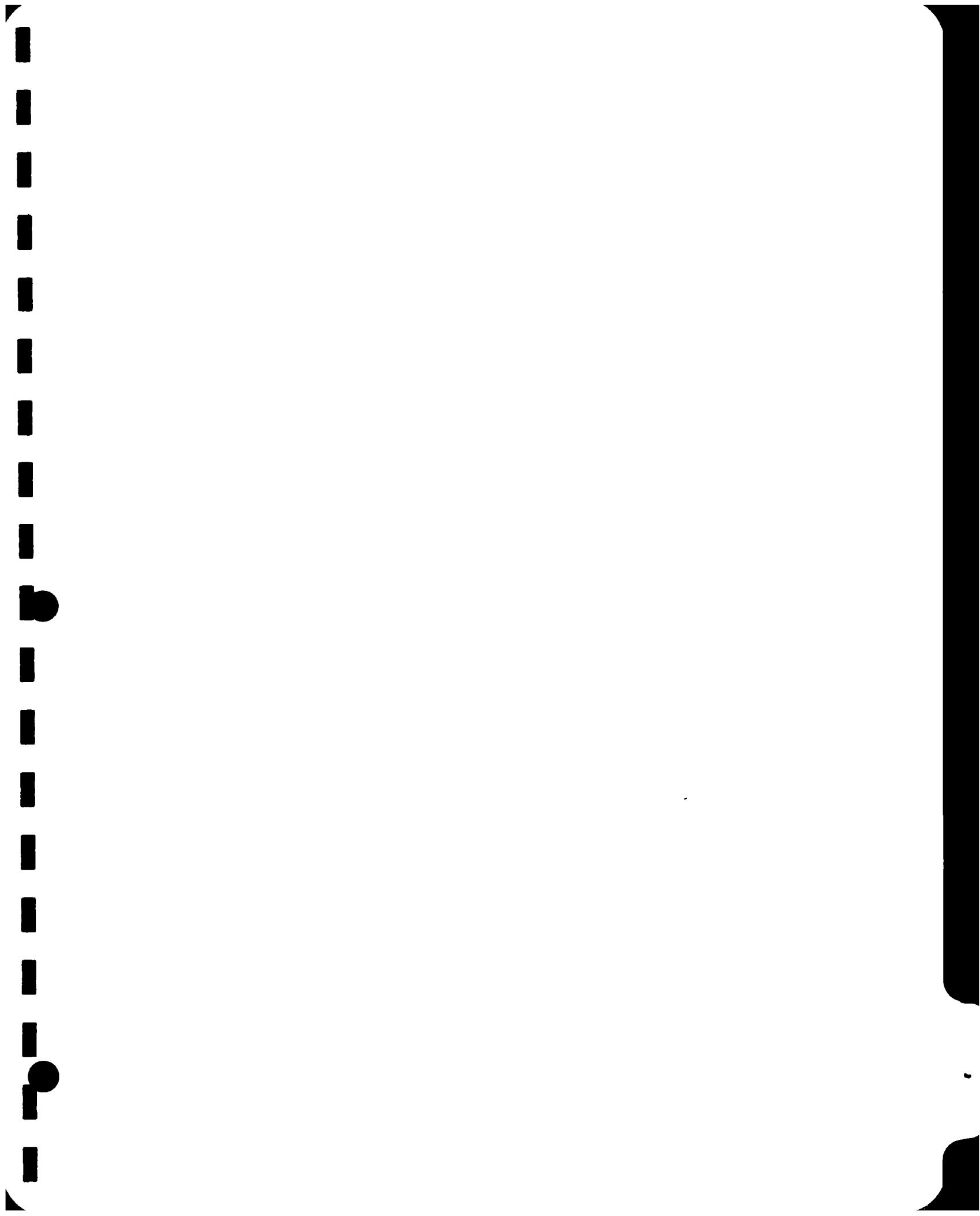
IMMEDIATE REMOVAL ACTION CHECK SHEET

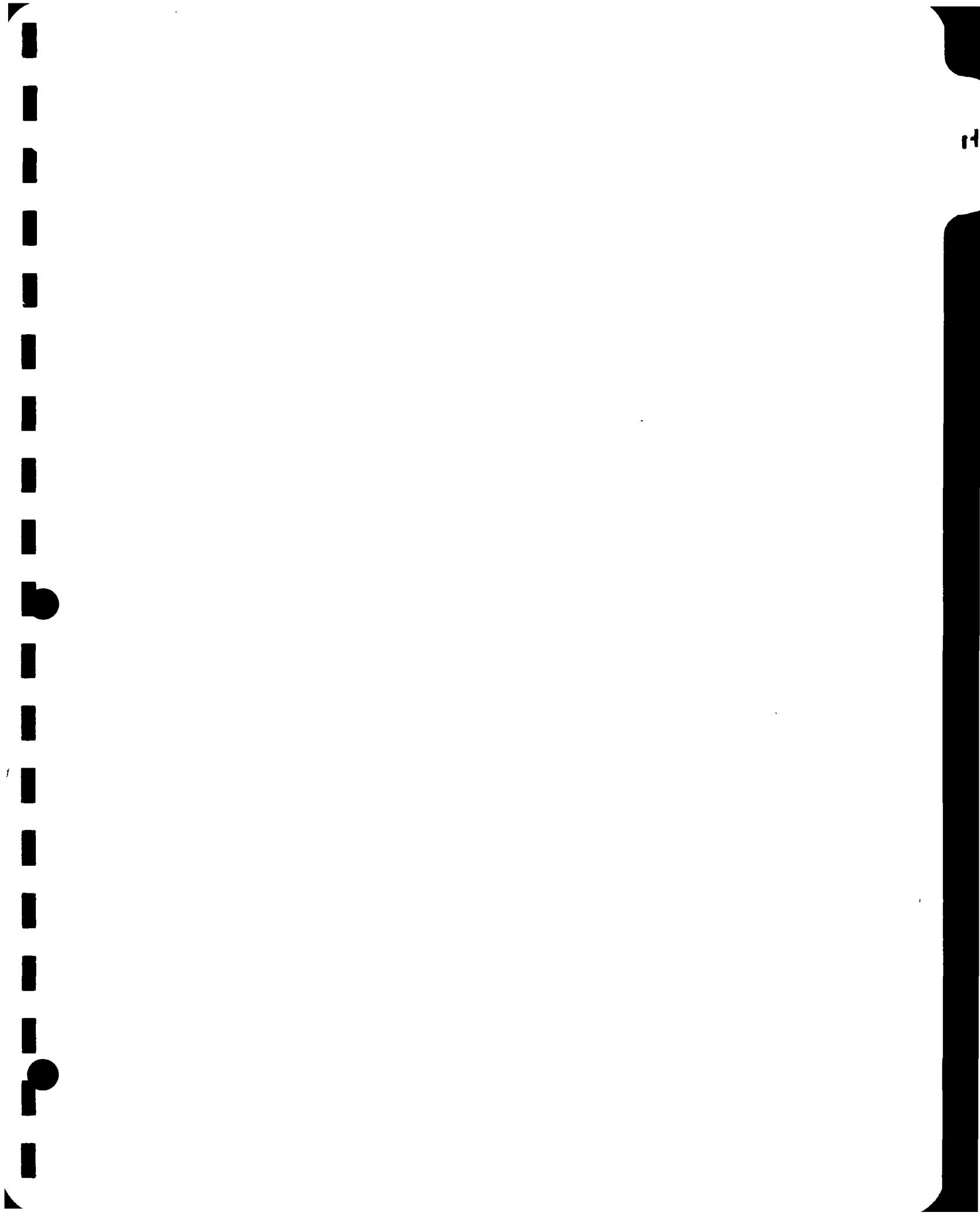
	High	Moderate	Low
<u>Fire and Explosion Hazard</u>			
Flammable Materials _____			X
Explosives _____			X
Incompatible Chemicals _____			X
<u>Direct Contact With Acutely Toxic Chemicals</u>			
Site Security _____			X
Leaking Drums or Tanks _____			X
Open Lagoons or Pits _____			X
Materials on Surface _____			X
Proximity of Population _____			X
Evidence of Casual Site Use _____			X
<u>Contaminated Water Supply</u>			
Exceeds 10 Day Snarl _____			X
Gross Taste or Odors _____	X		
Alternate Water Available _____			X
Potential Contamination _____			X
Is the site abandoned, active, or inactive? _____ <i>Inactive</i>			

Comments:

NFRAP

*SHP*





228320

184

2 different tillers

CJ. AND OBERT

ASL IN HADER NIP:

NIP ~~100~~

Dave! Longworthy

Tiller, Obert, Hader, Minn Mar 61

103 ft deep 6" well to limestone  
38 ft to water 35 ft to limestone  
Cased 45 ft with 6 in.  
Cased 10 ft into Galena 45' with 6" pipe  
pump set 53

8000 1-3.5 duct

0 GAL 35-103 Galena 7/155

11-16-31 A

SL 119°

70-A

neighbor  
70A

No. Tiller's in

Aquifer  
CAL-GAL

Ack Hader

**CODED**

open hole

Goodman  
non responsive

FORMATION	DEPTH	FORMATION	FROM	TO	
Quuu clay				24	
↓ shale				24	47
QTUU gravel & broken limerock			47	53	
OPAL limerock	7/1157		53	115	
OPPL shale	7/1095		115	217	
CSTR sandrock	7/993		217	321	
OPDC limerock	7/887		321	390	
III-16-31 BDCDDA					
EL - 1210 ± 5'					
Aquifer OPDC - OPDC	70-A				
Owner		open well			
<b>CODED</b>					
Use a second sheet if needed					
15. REMARKS, ELEVATION, SOURCE OF DATA, etc					
16. WATER WELL CONTRACTOR'S CERTIFICATION					
This well was drilled under my jurisdiction and this report is true to the best of my knowledge and belief					
Common Well Drilling <span style="float: right;">26240 License No.</span>					
Address: RR1 Box 266A Cannon Falls, Mn.					
Signed: <i>P. Otto</i> Date: <i>7/16/78</i> Authorized Representative					
Frank Otto <span style="float: right;">Name of Driller</span> Date <i>12/16/78</i>					
162712					

File date 9/11-16-34

9/11-16-34 CCR 64

Ex. 12-1

64F 218799

S. - 100

6f

• Section

HANSON, HAROLD 9/11-16-34-BBA  
Belle Creek Township  
Goodhue, Minn.

7-67  
all 1200

336' well  
261' cased 6"  
234 $\frac{1}{2}$ ' to water

LOG:

Quv	0-40	Drift DRFT	T/1161
CDCR	40-125	Shale SHLE	
OPVL	125-141	Platteville LMSN	
O FWD	141-150	Glenwood SHLE	
OSTP	150-263	St. Peter SNS	T/1051
OPDC	263-335	Shakopee DLNT	T/938

Hardness--19ppm; iron 10''m; pH 7.0 neutral

Aquifer  
OPDC - CDC

**CODED**

## Minnesota WELL LOG

**CANNON WELL COMPANY**

Cannon Falls, Minn. 55009

non responsive

Well Owner: \_\_\_\_\_

Location: \_\_\_\_\_

Date Completed: 12-14-73 Driller: Ray

Depth Description of Formation

0 to 25. dirt &amp; clay

25 to 220. sand rock

220 to 310. lime rock

to \_\_\_\_\_.

## SIZES AND MATERIALS USED

Top Casing Line: Inside Diameter 5 in. Wt. per ft. 15 lbs.  
Depth 225 ft. in.

Any Reduced Casing Sizes 5

Total Depth to bottom of Casing 81 ft. in.

Total Depth to bottom of Well 107 ft. in.

Depth measured from: ground Water level:

Screened Well: Size of Screen: Diam. in. Length ft. Slot \_\_\_\_\_

Make of Screen Metal

Fittings

Rock Well: Open Borehole 5 inches diam. ft. deep below casing

Test data: (gpm) ft. drawdown. Pumped for hrs.  
(gph)

Test Pump: \_\_\_\_\_

Tot. Len. of Setting: \_\_\_\_\_

NOTES: \_\_\_\_\_

## Buck Creek WELL LOG

**CANNON WELL COMPANY**

Cannon Falls, Minn. 55009

non responsive

Well Owner: \_\_\_\_\_

Location: \_\_\_\_\_

Date Completed: 6-12-71 Driller: John

Depth Description of Formation

0 to 81. old well

81 to 107. lime-rock or sandrock

to \_\_\_\_\_.

## SIZES AND MATERIALS USED

Top Casing Line: Inside Diameter \_\_\_\_\_ in. Wt. per ft. \_\_\_\_\_ lbs.  
Depth \_\_\_\_\_ ft. \_\_\_\_\_ in.

Any Reduced Casing Sizes 4

Total Depth to bottom of Casing 84 ft. in.

Total Depth to bottom of Well 107 ft. in.

Depth measured from: ground Water level:

Screened Well: Size of Screen: Diam. in. Length ft. Slot \_\_\_\_\_

Make of Screen Metal

Fittings

Rock Well: Open Borehole 4 inches diam. 25 ft. deep below casing

Test data: (gpm) ft. drawdown. Pumped for hrs.  
(gph)

Test Pump: \_\_\_\_\_

Tot. Len. of Setting: \_\_\_\_\_

NOTES: 4 Redrilled to try and get rid of oil-tast

still bad.

*Wells*  
*Minnola* WELL LOG

# CANNON WELL COMPANY

Cannon Falls, Minn. 55009

Well Owner: non responsive

Location:

Date Completed: 9-21-70 Driller: Bob Vaught

Depth	Description of Formation
0 to 13	Dirt & clay
13 to 15	Limestone
15 to 45	Blue clay
45 to 99	Layered limestone & blue clay
99 to 115	Limestone (very hard)
115 to 124	Gray clay
124 to 146	Gray sandrock
146 to 155	Tan sandrock
155 to 188	Yellow sandrock
188 to 210	Tan sandrock

#### SIZES AND MATERIALS USED

Top Casing Line: Inside Diameter 5 in. Wt. per ft. 15 lbs.  
Depth 245 ft. 6 in.

Any Reduced Casing Sizes --

Total Depth to bottom of Casing 245 ft. 6 in.

Total Depth to bottom of Well 260 ft. in.

Depth measured from: ground Water level: 180

Screened Well: Size of Screen: Diam. -- in. Length ft. Slot --

Make of Screen --- Metal ---

Fittings ---

Rock Well: Open Borehole 5 inches diam. 15 ft. deep below casing

Test data: 12 (gpm) ft. drawdown. Pumped for 1 hrs.  
(gph)

Test Pump: 3/4 HP submersible

Tot. Len. of Setting: 200 ft.

NOTES: \_\_\_\_\_

WELL LOG

# CANNON WELL COMPANY

Cannon Falls, Minn. 55009

Well Owner: non responsive

Location:

Date Completed: \_\_\_\_\_ Driller: \_\_\_\_\_

Depth	Description of Formation
0 to 210	on page 1
210 to 232	Light gray sandrock
232 to 238	Dark gray sandrock
238 to 239	Brown limestone
239 to 240	Gray sandrock
240 to 260	Limestone
to	
to	
to	
to	

**STATE OF MINNESOTA**      **DEPARTMENT OF HEALTH**

**WATER WELL RECORD**

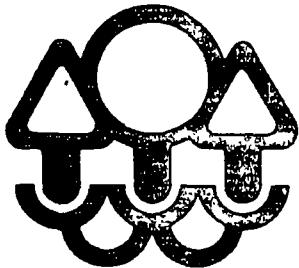
Minnesota Statute 166A.01-03

**MINNESOTA UNIQUE WELL NO**

### *for Water Samples*

435190

**non responsive**



## Minnesota Pollution Control Agency

December 29, 1988

Ms. Jeanne Griffin  
U.S. Environmental Protection Agency  
Waste Management Division  
Program Support Section (5HR-11)  
Remedial Response Branch  
230 South Dearborn Street  
Chicago, Illinois 60604

Dear Ms. Griffin:

Re: Minnesota Pollution Control Agency (MPCA) Submittal of Screening Site  
Inspection Reports  
Preliminary Assessment/Site Inspection (PA/SI) Cooperative Agreement with  
the U.S. Environmental Protection Agency

Pursuant to the PA/SI Cooperative Agreement, MPCA staff hereby submit for your  
review and approval Screening Site Inspection Reports for the following four  
( 4 ) potential hazardous waste sites:

<u>Site Name</u>	<u>EPA Identification Number</u>
1. FOLEY RR IMPROVEMENT PROJECT	MND981961659
2. HADER DUMP	MND981961865
3. HADER GROUND WATER CONTAMINATION	MND981961873
4. HOBART CORPORATION	MND068152768

If you have any comments or questions regarding these Screening Site Inspection  
Reports please contact me at 612/297-1793.

Sincerely,

Ronald R. Swenson  
Supervisor, Site Assessment Unit  
Program Development Section  
Ground Water and Solid Waste Division

RRS:

Enclosure

Phone: \_\_\_\_\_

520 Lafayette Road, St. Paul, Minnesota 55155  
Regional Offices • Duluth/Brainerd/Detroit Lakes/Marshall/Rochester  
Equal Opportunity Employer